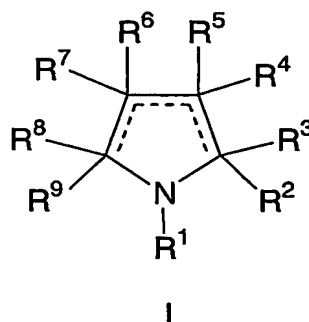


## WHAT IS CLAIMED IS:

1. A compound of Formula I:



- 5 or a pharmaceutically acceptable salt or stereoisomer thereof, wherein

- a is 0 or 1;  
 b is 0 or 1;  
 m is 0, 1, or 2;  
 10 n is 0 or 1;  
 r is 0 or 1;  
 s is 0 or 1;  
 u is 2, 3, 4 or 5;
- 15 a dashed line represents an optional double bond, provided that one and only one double bond is present in the ring;

R¹ is selected from:

- 1) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>1</sub>-C<sub>10</sub> alkyl,  
 20 2) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)aryl,  
 3) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>2</sub>-C<sub>10</sub> alkenyl,  
 4) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>2</sub>-C<sub>10</sub> alkynyl,  
 5) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>3</sub>-C<sub>8</sub> cycloalkyl,  
 6) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)heterocyclyl,  
 25 7) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)NR<sup>c</sup>R<sup>c'</sup>,  
 8) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>NR<sup>c</sup>R<sup>c'</sup>,  
 9) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,

- 10) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 11) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 12) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>-aryl,
- 13) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>-heterocyclyl,
- 5 14) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>-C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 15) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>P(=O)R<sup>d</sup>R<sup>d'</sup>,
- 16) aryl;
- 17) heterocyclyl; and
- 18) C<sub>1</sub>-C<sub>10</sub> alkyl;

10 said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, alkylene, heteroaryl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>2</sup> and R<sup>6</sup> are independently selected from:

- 1) aryl,
- 15 2) C<sub>1</sub>-C<sub>6</sub> aralkyl,
- 3) C<sub>3</sub>-C<sub>8</sub> cycloalkyl, and
- 4) heterocyclyl,

said aryl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

20

R<sup>3</sup>, R<sup>4</sup>, R<sup>5</sup>, R<sup>7</sup>, R<sup>8</sup>, and R<sup>9</sup> are independently selected from:

- 1) H,
- 2) C<sub>1</sub>-C<sub>10</sub> alkyl,
- 3) aryl,
- 25 4) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 5) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 7) C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,
- 8) C<sub>1</sub>-C<sub>6</sub> aralkyl,
- 9) C<sub>3</sub>-C<sub>8</sub> cycloalkyl, and
- 30 10) heterocyclyl,

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>; or

R<sup>4</sup> and R<sup>5</sup>, or R<sup>8</sup> and R<sup>9</sup>, attached to the same carbon atom are combined to form

$-(CH_2)_u-$  wherein one of the carbon atoms is optionally replaced by a moiety selected from O,  $S(O)_m$ ,  $-N(R^a)C(O)-$ ,  $-N(R^b)-$  and  $-N(COR^a)-$ ;

$R^{10}$  is independently selected from:

- |    |                                     |
|----|-------------------------------------|
| 5  | 1) $(C=O)_aO_bC_1-C_{10}$ alkyl,    |
|    | 2) $(C=O)_aO_b$ aryl,               |
|    | 3) $C_2-C_{10}$ alkenyl,            |
|    | 4) $C_2-C_{10}$ alkynyl,            |
|    | 5) $(C=O)_aO_b$ heterocyclyl,       |
| 10 | 6) $CO_2H$ ,                        |
|    | 7) halo,                            |
|    | 8) CN,                              |
|    | 9) OH,                              |
|    | 10) $O_bC_1-C_6$ perfluoroalkyl,    |
| 15 | 11) $O_a(C=O)_bNR^{12}R^{13}$ ,     |
|    | 12) $S(O)_mR^a$ ,                   |
|    | 13) $S(O)_2NR^{12}R^{13}$ ,         |
|    | 14) oxo,                            |
|    | 15) CHO,                            |
| 20 | 16) $(N=O)R^{12}R^{13}$ , or        |
|    | 17) $(C=O)_aO_bC_3-C_8$ cycloalkyl, |

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one or more substituents selected from  $R^{11}$ ;

25  $R^{11}$  is selected from:

- |    |  |
|----|--|
| 25 | 1) $(C=O)_rO_s(C_1-C_{10})$ alkyl,     |
|    | 2) $O_r(C_1-C_3)$ perfluoroalkyl,      |
|    | 3) $(C_0-C_6)$ alkylene- $S(O)_mR^a$ , |
|    | 4) oxo,                                |
| 30 | 5) OH,                                 |
|    | 6) halo,                               |
|    | 7) CN,                                 |
|    | 8) $(C=O)_rO_s(C_2-C_{10})$ alkenyl,   |
|    | 9) $(C=O)_rO_s(C_2-C_{10})$ alkynyl,   |

- 10)  $(C=O)_rO_s(C_3-C_6)cycloalkyl$ ,
- 11)  $(C=O)_rO_s(C_0-C_6)alkylene-aryl$ ,
- 12)  $(C=O)_rO_s(C_0-C_6)alkylene-heterocyclyl$ ,
- 13)  $(C=O)_rO_s(C_0-C_6)alkylene-N(R^b)_2$ ,
- 5 14)  $C(O)R^a$ ,
- 15)  $(C_0-C_6)alkylene-CO_2R^a$ ,
- 16)  $C(O)H$ ,
- 17)  $(C_0-C_6)alkylene-CO_2H$ ,
- 18)  $C(O)N(R^b)_2$ ,
- 10 19)  $S(O)_mR^a$ , and
- 20)  $S(O)_2N(R^b)_2$

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, alkylene and heterocyclyl is optionally substituted with up to three substituents selected from  $R^b$ , OH,  $(C_1-C_6)alkoxy$ , halogen,  $CO_2H$ , CN,  $O(C=O)C_1-C_6$  alkyl, oxo, and  $N(R^b)_2$ ;

15

$R^{12}$  and  $R^{13}$  are independently selected from:

- 1) H,
- 2)  $(C=O)O_bC_1-C_{10}$  alkyl,
- 3)  $(C=O)O_bC_3-C_8$  cycloalkyl,
- 20 4)  $(C=O)O_baryl$ ,
- 5)  $(C=O)O_bheterocyclyl$ ,
- 6)  $C_1-C_{10}$  alkyl,
- 7) aryl,
- 8)  $C_2-C_{10}$  alkenyl,
- 25 9)  $C_2-C_{10}$  alkynyl,
- 10) heterocyclyl,
- 11)  $C_3-C_8$  cycloalkyl,
- 12)  $SO_2R^a$ , and
- 13)  $(C=O)NR^b_2$ ,

- 30 said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one or more substituents selected from  $R^{11}$ , or

$R^{12}$  and  $R^{13}$  can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and

optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one or more substituents selected from R<sup>11</sup>;

5 R<sup>14</sup> is independently selected from:

- 1) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 2) (C=O)<sub>a</sub>O<sub>b</sub>aryl,
- 3) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 4) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 10 5) (C=O)<sub>a</sub>O<sub>b</sub> heterocyclyl,
- 6) CO<sub>2</sub>H,
- 7) halo,
- 8) CN,
- 9) OH,
- 15 10) O<sub>b</sub>C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,
- 11) O<sub>a</sub>(C=O)<sub>b</sub>NR<sup>12</sup>R<sup>13</sup>,
- 12) S(O)<sub>m</sub>R<sup>a</sup>,
- 13) S(O)<sub>2</sub>NR<sup>12</sup>R<sup>13</sup>,
- 14) oxo,
- 20 15) CHO,
- 16) (N=O)R<sup>12</sup>R<sup>13</sup>, or
- 17) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one or more substituents selected from R<sup>11</sup>;

25

R<sup>a</sup> is (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, aryl, or heterocyclyl, optionally substituted with one to three substituents selected from R<sup>14</sup>;

30 R<sup>b</sup> is H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, (C=O)OC<sub>1</sub>-C<sub>6</sub> alkyl, (C=O)C<sub>1</sub>-C<sub>6</sub> alkyl or S(O)<sub>2</sub>R<sup>a</sup>, optionally substituted with one to three substituents selected from R<sup>14</sup>;

$R^c$  and  $R^{c'}$  are independently selected from: H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl and (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, optionally substituted with one, two or three substituents selected from  $R^{10}$ , or

$R^c$  and  $R^{c'}$  can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from  $R^{11}$ ;

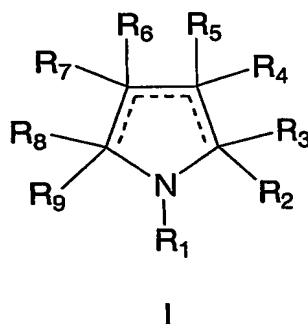
$R^d$  and  $R^{d'}$  are independently selected from: (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>1</sub>-C<sub>6</sub>)alkoxy and  $NR^{b2}$ , or

$R^d$  and  $R^{d'}$  can be taken together with the phosphorous to which they are attached to form a monocyclic heterocycle with 5-7 members the ring and optionally containing, in addition to the phosphorous, one or two additional heteroatoms selected from  $NR^e$ , O and S, said monocyclic heterocycle optionally substituted with one, two or three substituents selected from  $R^{11}$ ;

$R^e$  is selected from: H and (C<sub>1</sub>-C<sub>6</sub>)alkyl; and

X is selected from O,  $NR^e$  and S.

2. The compound according to Claim 1 of the Formula I:



or a pharmaceutically acceptable salt or stereoisomer thereof, wherein

a is 0 or 1;

- b is 0 or 1;  
 m is 0, 1, or 2;  
 n is 0 or 1;  
 r is 0 or 1;  
 5 s is 0 or 1;  
 u is 2, 3, 4 or 5;

a dashed line represents an optional double bond, provided that one and only one double bond is present in the ring;

10

R<sup>1</sup> is selected from:

- 1) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>1</sub>-C<sub>10</sub> alkyl,
  - 2) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)aryl,
  - 3) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>2</sub>-C<sub>10</sub> alkenyl,
  - 15 4) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>2</sub>-C<sub>10</sub> alkynyl,
  - 5) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
  - 6) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)heterocyclyl,
  - 7) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)NR<sup>c</sup>R<sup>c'</sup>,
  - 8) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>NR<sup>c</sup>R<sup>c'</sup>,
  - 20 9) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
  - 10) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>C<sub>2</sub>-C<sub>10</sub> alkenyl,
  - 11) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>C<sub>2</sub>-C<sub>10</sub> alkynyl,
  - 12) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>-aryl,
  - 13) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>-heterocyclyl,
  - 25 14) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>-C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
  - 15) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>P(=O)R<sup>d</sup>R<sup>d'</sup>,
  - 16) aryl;
  - 17) heterocyclyl; and
  - 18) C<sub>1</sub>-C<sub>10</sub> alkyl;
- 30 said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, alkylene, heteroaryl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>2</sup> and R<sup>6</sup> are independently selected from:

- 1) aryl,
- 35 2) C<sub>1</sub>-C<sub>6</sub> aralkyl,

3) C<sub>3</sub>-C<sub>8</sub> cycloalkyl, and

4) heterocyclyl,

said aryl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

5

R<sup>3</sup>, R<sup>4</sup>, R<sup>5</sup>, R<sup>7</sup>, R<sup>8</sup>, and R<sup>9</sup> are independently selected from:

1) H,

2) C<sub>1</sub>-C<sub>10</sub> alkyl,

3) aryl,

10 4) C<sub>2</sub>-C<sub>10</sub> alkenyl,

5) C<sub>2</sub>-C<sub>10</sub> alkynyl,

6) C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,

7) C<sub>1</sub>-C<sub>6</sub> aralkyl,

8) C<sub>3</sub>-C<sub>8</sub> cycloalkyl, and

15 9) heterocyclyl,

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>; or

20 R<sup>4</sup> and R<sup>5</sup>, or R<sup>8</sup> and R<sup>9</sup>, attached to the same carbon atom are combined to form -(CH<sub>2</sub>)<sub>u</sub>- wherein one of the carbon atoms is optionally replaced by a moiety selected from O, S(O)<sub>m</sub>, -N(R<sup>a</sup>)C(O)-, -N(R<sup>b</sup>)- and -N(COR<sup>a</sup>)-;

R<sup>10</sup> is independently selected from:

1) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,

25 2) (C=O)<sub>a</sub>O<sub>b</sub>aryl,

3) C<sub>2</sub>-C<sub>10</sub> alkenyl,

4) C<sub>2</sub>-C<sub>10</sub> alkynyl,

5) (C=O)<sub>a</sub>O<sub>b</sub> heterocyclyl,

6) CO<sub>2</sub>H,

30 7) halo,

8) CN,

9) OH,

10) O<sub>b</sub>C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,

11) O<sub>a</sub>(C=O)<sub>b</sub>NR<sup>12</sup>R<sup>13</sup>,



- 12)  $S(O)_m R^a$ ,  
 13)  $S(O)_2 N R^{12} R^{13}$ ,  
 14) oxo,  
 15) CHO,  
 5 16)  $(N=O) R^{12} R^{13}$ , or  
 17)  $(C=O)_a O_b C_3-C_8$  cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one or more substituents selected from  $R^{11}$ ;

10  $R^{11}$  is selected from:

- 1)  $(C=O)_r O_s (C_1-C_{10})$  alkyl,  
 2)  $O_r (C_1-C_3)$  perfluoroalkyl,  
 3)  $(C_0-C_6)$  alkylene- $S(O)_m R^a$ ,  
 4) oxo,  
 15 5) OH,  
 6) halo,  
 7) CN,  
 8)  $(C=O)_r O_s (C_2-C_{10})$  alkenyl,  
 9)  $(C=O)_r O_s (C_2-C_{10})$  alkynyl,  
 20 10)  $(C=O)_r O_s (C_3-C_6)$  cycloalkyl,  
 11)  $(C=O)_r O_s (C_0-C_6)$  alkylene-aryl,  
 12)  $(C=O)_r O_s (C_0-C_6)$  alkylene-heterocyclyl,  
 13)  $(C=O)_r O_s (C_0-C_6)$  alkylene- $N(R^b)_2$ ,  
 14)  $C(O) R^a$ ,  
 25 15)  $(C_0-C_6)$  alkylene- $CO_2 R^a$ ,  
 16)  $C(O)H$ ,  
 17)  $(C_0-C_6)$  alkylene- $CO_2H$ ,  
 18)  $C(O)N(R^b)_2$ ,  
 19)  $S(O)_m R^a$ , and  
 30 20)  $S(O)_2 N(R^b)_2$

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, alkylene and heterocyclyl is optionally substituted with up to three substituents selected from  $R^b$ , OH,  $(C_1-C_6)$  alkoxy, halogen,  $CO_2H$ , CN,  $O(C=O)C_1-C_6$  alkyl, oxo, and  $N(R^b)_2$ ;

R<sup>12</sup> and R<sup>13</sup> are independently selected from:

- 1) H,
  - 2) (C=O)O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
  - 3) (C=O)O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
  - 5 4) (C=O)O<sub>b</sub>aryl,
  - 5) (C=O)O<sub>b</sub>heterocyclyl,
  - 6) C<sub>1</sub>-C<sub>10</sub> alkyl,
  - 7) aryl,
  - 8) C<sub>2</sub>-C<sub>10</sub> alkenyl,
  - 10 9) C<sub>2</sub>-C<sub>10</sub> alkynyl,
  - 10) heterocyclyl,
  - 11) C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
  - 12) SO<sub>2</sub>R<sup>a</sup>, and
  - 13) (C=O)NR<sup>b</sup><sub>2</sub>,
- 15 said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one or more substituents selected from R<sup>11</sup>, or

R<sup>12</sup> and R<sup>13</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and  
 20 optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one or more substituents selected from R<sup>11</sup>;

R<sup>a</sup> is (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, aryl, or heterocyclyl;

25 R<sup>b</sup> is H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, (C=O)OC<sub>1</sub>-C<sub>6</sub> alkyl, (C=O)C<sub>1</sub>-C<sub>6</sub> alkyl or S(O)<sub>2</sub>R<sup>a</sup>;

R<sup>c</sup> and R<sup>c'</sup> are independently selected from: H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl and  
 30 (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, optionally substituted with one, two or three substituents selected from R<sup>10</sup>, or

R<sup>c</sup> and R<sup>c'</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally

containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

- 5 R<sup>d</sup> and R<sup>d'</sup> are independently selected from: (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>1</sub>-C<sub>6</sub>)alkoxy and NR<sup>b</sup><sub>2</sub>, or

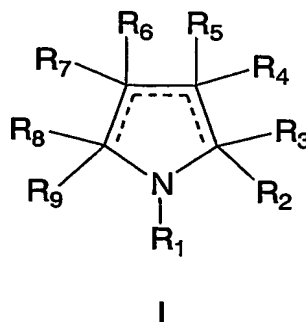
- R<sup>d</sup> and R<sup>d'</sup> can be taken together with the phosphorous to which they are attached to form a monocyclic heterocycle with 5-7 members the ring and optionally containing,  
 10 in addition to the phosphorous, one or two additional heteroatoms selected from NR<sup>e</sup>, O and S, said monocyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

R<sup>e</sup> is selected from: H and (C<sub>1</sub>-C<sub>6</sub>)alkyl; and

15

X is selected from O, NR<sup>e</sup> and S.

3. The compound according to Claim 2 of Formula I:



- 20 or a pharmaceutically acceptable salt or stereoisomer thereof, wherein

- a is 0 or 1;  
 b is 0 or 1;  
 m is 0, 1, or 2;  
 25 n is 0 or 1;  
 r is 0 or 1;  
 s is 0 or 1;

u is 2, 3, 4 or 5;

a dashed line represents an optional double bond, provided that one and only one double bond is present in the ring;

5

R<sup>1</sup> is selected from:

- 1) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>1</sub>-C<sub>10</sub> alkyl,
- 2) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)aryl,
- 3) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 10 4) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 5) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 6) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)heterocyclyl,
- 7) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)NR<sup>c</sup>R<sup>c'</sup>,
- 8) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>NR<sup>c</sup>R<sup>c'</sup>,
- 15 9) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 10) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 11) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 12) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>-aryl,
- 13) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>-heterocyclyl,
- 20 14) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>-C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 15) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>P(=O)R<sup>d</sup>R<sup>d'</sup>,
- 16) aryl;
- 17) heterocyclyl; and
- 18) C<sub>1</sub>-C<sub>10</sub> alkyl;

25 said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, alkylene, heteroaryl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>2</sup> and R<sup>6</sup> are independently selected from:

- 1) aryl,
- 30 2) C<sub>1</sub>-C<sub>6</sub> aralkyl,
- 3) C<sub>3</sub>-C<sub>8</sub> cycloalkyl, and
- 4) heterocyclyl,

said aryl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

35

R<sup>3</sup>, R<sup>4</sup>, R<sup>5</sup>, R<sup>7</sup>, R<sup>8</sup>, and R<sup>9</sup> are independently selected from:

- 1) H,
- 2) C<sub>1</sub>-C<sub>10</sub> alkyl,
- 3) aryl,
- 5 4) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 5) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 6) C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,
- 7) C<sub>1</sub>-C<sub>6</sub> aralkyl,
- 8) C<sub>3</sub>-C<sub>8</sub> cycloalkyl, and
- 10 9) heterocyclyl,

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>; or

15 R<sup>4</sup> and R<sup>5</sup>, or R<sup>8</sup> and R<sup>9</sup>, attached to the same carbon atom are combined to form -(CH<sub>2</sub>)<sub>u</sub>- wherein one of the carbon atoms is optionally replaced by a moiety selected from O, S(O)<sub>m</sub>, -N(R<sup>a</sup>)C(O)-, -N(R<sup>b</sup>)- and -N(COR<sup>a</sup>)-;

R<sup>10</sup> is independently selected from:

- 1) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 2) (C=O)<sub>a</sub>O<sub>b</sub>aryl,
- 20 3) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 4) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 5) (C=O)<sub>a</sub>O<sub>b</sub> heterocyclyl,
- 6) CO<sub>2</sub>H,
- 7) halo,
- 25 8) CN,
- 9) OH,
- 10) O<sub>b</sub>C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,
- 11) O<sub>a</sub>(C=O)<sub>b</sub>NR<sup>12</sup>R<sup>13</sup>,
- 12) S(O)<sub>m</sub>R<sup>a</sup>,
- 30 13) S(O)<sub>2</sub>NR<sup>12</sup>R<sup>13</sup>,
- 14) oxo,
- 15) CHO,
- 16) (N=O)R<sup>12</sup>R<sup>13</sup>, or
- 17) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one or more substituents selected from R<sup>11</sup>;

R<sup>11</sup> is selected from:

- 5           1) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>1</sub>-C<sub>10</sub>)alkyl,
- 2) O<sub>r</sub>(C<sub>1</sub>-C<sub>3</sub>)perfluoroalkyl,
- 3) (C<sub>0</sub>-C<sub>6</sub>)alkylene-S(O)<sub>m</sub>R<sup>a</sup>,
- 4) oxo,
- 5) OH,
- 10          6) halo,
- 7) CN,
- 8) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>2</sub>-C<sub>10</sub>)alkenyl,
- 9) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>2</sub>-C<sub>10</sub>)alkynyl,
- 10) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
- 15          11) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-aryl,
- 12) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-heterocyclyl,
- 13) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-N(R<sup>b</sup>)<sub>2</sub>,
- 14) C(O)R<sup>a</sup>,
- 15) (C<sub>0</sub>-C<sub>6</sub>)alkylene-CO<sub>2</sub>R<sup>a</sup>,
- 20          16) C(O)H,
- 17) (C<sub>0</sub>-C<sub>6</sub>)alkylene-CO<sub>2</sub>H,
- 18) C(O)N(R<sup>b</sup>)<sub>2</sub>,
- 19) S(O)<sub>m</sub>R<sup>a</sup>, and
- 20) S(O)<sub>2</sub>N(R<sup>b</sup>)<sub>2</sub>,

- 25   said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, alkylene and heterocyclyl is optionally substituted with up to three substituents selected from R<sup>b</sup>, OH, (C<sub>1</sub>-C<sub>6</sub>)alkoxy, halogen, CO<sub>2</sub>H, CN, O(C=O)C<sub>1</sub>-C<sub>6</sub> alkyl, oxo, and N(R<sup>b</sup>)<sub>2</sub>;

R<sup>12</sup> and R<sup>13</sup> are independently selected from:

- 30          1) H,
- 2) (C=O)O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 3) (C=O)O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 4) (C=O)O<sub>b</sub>aryl,
- 5) (C=O)O<sub>b</sub>heterocyclyl,

- 5           6)     C<sub>1</sub>-C<sub>10</sub> alkyl,  
              7)     aryl,  
              8)     C<sub>2</sub>-C<sub>10</sub> alkenyl,  
              9)     C<sub>2</sub>-C<sub>10</sub> alkynyl,  
              10)    heterocyclyl,  
              11)    C<sub>3</sub>-C<sub>8</sub> cycloalkyl,  
              12)    SO<sub>2</sub>R<sup>a</sup>, and  
              13)    (C=O)NR<sup>b</sup><sub>2</sub>,

10   said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one or more substituents selected from R<sup>11</sup>, or

15   R<sup>12</sup> and R<sup>13</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one or more substituents selected from R<sup>11</sup>;

R<sup>a</sup> is (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, aryl, or heterocyclyl;

20   R<sup>b</sup> is H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, (C=O)OC<sub>1</sub>-C<sub>6</sub> alkyl, (C=O)C<sub>1</sub>-C<sub>6</sub> alkyl or S(O)<sub>2</sub>R<sup>a</sup>;

25   R<sup>c</sup> and R<sup>c'</sup> are independently selected from: H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl and (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl or

30   R<sup>c</sup> and R<sup>c'</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

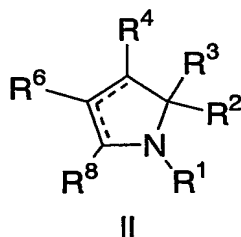
R<sup>d</sup> and R<sup>d'</sup> are independently selected from: (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>1</sub>-C<sub>6</sub>)alkoxy and NR<sup>b</sup><sub>2</sub>, or

$R^d$  and  $R^{d'}$  can be taken together with the phosphorous to which they are attached to form a monocyclic heterocycle with 5-7 members the ring and optionally containing, in addition to the phosphorous, one or two additional heteroatoms selected from  $NR^e$ , O and S, said monocyclic heterocycle optionally substituted with one, two or three substituents selected from  $R^{11}$ ;

$R^e$  is selected from: H and  $(C_1-C_6)$ alkyl; and

X is selected from O,  $NR^e$  and S.

4. The compound according to Claim 2 of the Formula II,



or a pharmaceutically acceptable salt or stereoisomer thereof, wherein

wherein:

a is 0 or 1;

b is 0 or 1;

m is 0, 1, or 2;

r is 0 or 1;

s is 0 or 1;

a dashed line represents an optional double bond, provided that one and only one double bond is present in the ring;

$R^1$  is selected from:

- 1)  $(C_1-C_6\text{-alkylene})_n(C=O)C_1-C_{10}$  alkyl,
- 2)  $(C_1-C_6\text{-alkylene})_n(C=O)$ aryl,
- 3)  $(C_1-C_6\text{-alkylene})_n(C=O)C_2-C_{10}$  alkenyl,



- 4) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=O)C<sub>2</sub>-C<sub>10</sub> alkynyl,  
 5) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=O)C<sub>3</sub>-C<sub>8</sub> cycloalkyl,  
 6) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=O)heterocyclyl,  
 7) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=O)NR<sup>c</sup>R<sup>c'</sup>,  
 5 8) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>NR<sup>c</sup>R<sup>c'</sup>,  
 9) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,  
 10) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>-aryl,  
 11) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>-heterocyclyl,  
 12) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>-C<sub>3</sub>-C<sub>8</sub> cycloalkyl,  
 10 13) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>P(=O)R<sup>d</sup>R<sup>d'</sup>,  
 14) aryl,  
 15) heterocyclyl, and  
 16) C<sub>1</sub>-C<sub>10</sub> alkyl;

15 said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, alkylene, heteroaryl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>2</sup> and R<sup>6</sup> are independently selected from:

- 1) aryl,  
 2) C<sub>1</sub>-C<sub>6</sub> aralkyl,  
 20 3) C<sub>3</sub>-C<sub>8</sub> cycloalkyl, and  
 4) heterocyclyl,

said aryl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

25 R<sup>3</sup>, R<sup>4</sup> and R<sup>8</sup> are independently selected from:

- 1) H,  
 2) C<sub>1</sub>-C<sub>10</sub> alkyl,  
 3) aryl,  
 4) C<sub>2</sub>-C<sub>10</sub> alkenyl,  
 30 5) C<sub>2</sub>-C<sub>10</sub> alkynyl,  
 6) C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,  
 7) C<sub>1</sub>-C<sub>6</sub> aralkyl,  
 8) C<sub>3</sub>-C<sub>8</sub> cycloalkyl, and  
 9) heterocyclyl,

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>10</sup> is independently selected from:

- 5           1) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 2) (C=O)<sub>a</sub>O<sub>b</sub>aryl,
- 3) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 4) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 5) (C=O)<sub>a</sub>O<sub>b</sub> heterocyclyl,
- 10          6) CO<sub>2</sub>H,
- 7) halo,
- 8) CN,
- 9) OH,
- 10) O<sub>b</sub>C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,
- 15          11) O<sub>a</sub>(C=O)<sub>b</sub>NR<sup>12</sup>R<sup>13</sup>,
- 12) S(O)<sub>m</sub>R<sup>a</sup>,
- 13) S(O)<sub>2</sub>NR<sup>12</sup>R<sup>13</sup>,
- 14) oxo,
- 15) CHO,
- 20          16) (N=O)R<sup>12</sup>R<sup>13</sup>, or
- 17) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

25 R<sup>11</sup> is selected from:

- 1) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>1</sub>-C<sub>10</sub>)alkyl,
- 2) O<sub>r</sub>(C<sub>1</sub>-C<sub>3</sub>)perfluoroalkyl,
- 3) oxo,
- 4) OH,
- 30          5) halo,
- 6) CN,
- 7) (C<sub>2</sub>-C<sub>10</sub>)alkenyl,
- 8) (C<sub>2</sub>-C<sub>10</sub>)alkynyl,
- 9) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
- 35          10) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-aryl,

- 11)  $(\text{C}=\text{O})_r\text{O}_s(\text{C}_0\text{-C}_6)\text{alkylene-heterocyclyl}$ ,
- 12)  $(\text{C}=\text{O})_r\text{O}_s(\text{C}_0\text{-C}_6)\text{alkylene-N(R}^b)_2$ ,
- 13)  $\text{C(O)R}^a$ ,
- 14)  $(\text{C}_0\text{-C}_6)\text{alkylene-CO}_2\text{R}^a$ ,
- 5 15)  $\text{C(O)H}$ ,
- 16)  $(\text{C}_0\text{-C}_6)\text{alkylene-CO}_2\text{H}$ ,
- 17)  $\text{C(O)N(R}^b)_2$ ,
- 18)  $\text{S(O)}_m\text{R}^a$ , and
- 19)  $\text{S(O)}_2\text{N(R}^b)_2$ ;

10

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, alkylene and heterocyclyl is optionally substituted with up to three substituents selected from  $\text{R}^b$ , OH,  $(\text{C}_1\text{-C}_6)\text{alkoxy}$ , halogen,  $\text{CO}_2\text{H}$ , CN,  $\text{O}(\text{C}=\text{O})\text{C}_1\text{-C}_6$  alkyl, oxo, and  $\text{N(R}^b)_2$ ;

15  $\text{R}^{12}$  and  $\text{R}^{13}$  are independently selected from:

- 1) H,
- 2)  $(\text{C}=\text{O})\text{O}_b\text{C}_1\text{-C}_{10}$  alkyl,
- 3)  $(\text{C}=\text{O})\text{O}_b\text{C}_3\text{-C}_8$  cycloalkyl,
- 4)  $(\text{C}=\text{O})\text{O}_b\text{aryl}$ ,
- 20 5)  $(\text{C}=\text{O})\text{O}_b\text{heterocyclyl}$ ,
- 6)  $\text{C}_1\text{-C}_{10}$  alkyl,
- 7) aryl,
- 8)  $\text{C}_2\text{-C}_{10}$  alkenyl,
- 9)  $\text{C}_2\text{-C}_{10}$  alkynyl,
- 25 10) heterocyclyl,
- 11)  $\text{C}_3\text{-C}_8$  cycloalkyl,
- 12)  $\text{SO}_2\text{R}^a$ , and
- 13)  $(\text{C}=\text{O})\text{NR}^b_2$ ,

30 said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one, two or three substituents selected from  $\text{R}^{11}$ , or

$\text{R}^{12}$  and  $\text{R}^{13}$  can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms

selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

R<sup>a</sup> is (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, aryl, or heterocyclyl;

5

R<sup>b</sup> is H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, (C=O)OC<sub>1</sub>-C<sub>6</sub> alkyl, (C=O)C<sub>1</sub>-C<sub>6</sub> alkyl or S(O)<sub>2</sub>R<sup>a</sup>;

10 R<sup>c</sup> and R<sup>c'</sup> are independently selected from: H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl and (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl; or R<sup>c</sup> and R<sup>c'</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

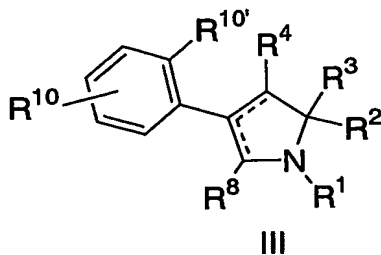
15

R<sup>d</sup> and R<sup>d'</sup> are independently selected from: (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>1</sub>-C<sub>6</sub>)alkoxy and NR<sup>b</sup><sub>2</sub>, or

20 R<sup>d</sup> and R<sup>d'</sup> can be taken together with the phosphorous to which they are attached to form a monocyclic heterocycle with 5-7 members the ring and optionally containing, in addition to the phosphorous, one or two additional heteroatoms selected from NRe, O and S, said monocyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>; and

25 R<sup>e</sup> is selected from: H and (C<sub>1</sub>-C<sub>6</sub>)alkyl.

5. A compound of the Formula III,



or a pharmaceutically acceptable salt or stereoisomer thereof, wherein

- a is 0 or 1;  
 b is 0 or 1;  
 m is 0, 1, or 2;  
 5 r is 0 or 1;  
 s is 0 or 1;

a dashed line represents an optional double bond, provided that one and only one double bond is present in the ring;

10

R<sup>1</sup> is selected from:

- 1) (C=O)C<sub>1</sub>-C<sub>10</sub> alkyl,
- 2) (C=O)aryl,
- 3) (C=O)C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 15 4) (C=O)C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 5) (C=O)C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 6) (C=O)heterocyclyl,
- 7) (C=O)NR<sup>c</sup>R<sup>c'</sup>,
- 8) SO<sub>2</sub>NR<sup>c</sup>R<sup>c'</sup>,
- 20 9) SO<sub>2</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 10) SO<sub>2</sub>-aryl,
- 11) SO<sub>2</sub>-heterocyclyl,
- 12) SO<sub>2</sub>-C<sub>3</sub>-C<sub>8</sub> cycloalkyl, and
- 13) P(=O)R<sup>d</sup>R<sup>d'</sup>,

- 25 said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, heteroaryl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>2</sup> is selected from:

- 1) aryl,
- 30 2) C<sub>1</sub>-C<sub>6</sub> aralkyl,
- 3) C<sub>3</sub>-C<sub>8</sub> cycloalkyl, and
- 4) heterocyclyl,

said aryl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

- 35 R<sup>3</sup>, R<sup>4</sup> and R<sup>8</sup> are independently selected from:

- 1) H,
  - 2) C<sub>1</sub>-C<sub>10</sub> alkyl,
  - 3) aryl,
  - 4) C<sub>2</sub>-C<sub>10</sub> alkenyl,
  - 5) C<sub>2</sub>-C<sub>10</sub> alkynyl,
  - 6) C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,
  - 7) C<sub>1</sub>-C<sub>6</sub> aralkyl,
  - 8) C<sub>3</sub>-C<sub>8</sub> cycloalkyl, and
  - 9) heterocyclyl,
- 10 said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>10</sup> is independently selected from:

- 1) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 15 2) (C=O)<sub>a</sub>O<sub>b</sub>aryl,
- 3) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 4) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 5) (C=O)<sub>a</sub>O<sub>b</sub> heterocyclyl,
- 6) CO<sub>2</sub>H,
- 20 7) halo,
- 8) CN,
- 9) OH,
- 10) O<sub>b</sub>C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,
- 11) O<sub>a</sub>(C=O)<sub>b</sub>NR<sup>12</sup>R<sup>13</sup>,
- 25 12) S(O)<sub>m</sub>R<sup>a</sup>,
- 13) S(O)<sub>2</sub>NR<sup>12</sup>R<sup>13</sup>,
- 14) oxo,
- 15) CHO,
- 16) (N=O)R<sup>12</sup>R<sup>13</sup>, or
- 30 17) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

R<sup>10'</sup> is halogen;

35

R<sup>11</sup> is selected from:

- 1) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>1</sub>-C<sub>10</sub>)alkyl,
- 2) O<sub>r</sub>(C<sub>1</sub>-C<sub>3</sub>)perfluoroalkyl,
- 3) oxo,
- 5 4) OH,
- 5) halo,
- 6) CN,
- 7) (C<sub>2</sub>-C<sub>10</sub>)alkenyl,
- 8) (C<sub>2</sub>-C<sub>10</sub>)alkynyl,
- 10 9) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
- 10) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-aryl,
- 11) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-heterocyclyl,
- 12) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-N(R<sup>b</sup>)<sub>2</sub>,
- 13) C(O)R<sup>a</sup>,
- 15 14) (C<sub>0</sub>-C<sub>6</sub>)alkylene-CO<sub>2</sub>R<sup>a</sup>,
- 15) C(O)H,
- 16) (C<sub>0</sub>-C<sub>6</sub>)alkylene-CO<sub>2</sub>H, and
- 17) C(O)N(R<sup>b</sup>)<sub>2</sub>,
- 18) S(O)<sub>m</sub>R<sup>a</sup>, and
- 20 19) S(O)<sub>2</sub>N(R<sup>b</sup>)<sub>2</sub>;

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, and heterocyclyl is optionally substituted with up to three substituents selected from R<sup>b</sup>, OH, (C<sub>1</sub>-C<sub>6</sub>)alkoxy, halogen, CO<sub>2</sub>H, CN, O(C=O)C<sub>1</sub>-C<sub>6</sub> alkyl, oxo, and N(R<sup>b</sup>)<sub>2</sub>;

25

R<sup>12</sup> and R<sup>13</sup> are independently selected from:

- 1) H,
- 2) (C=O)O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 3) (C=O)O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 30 4) (C=O)O<sub>b</sub>aryl,
- 5) (C=O)O<sub>b</sub>heterocyclyl,
- 6) C<sub>1</sub>-C<sub>10</sub> alkyl,
- 7) aryl,
- 8) C<sub>2</sub>-C<sub>10</sub> alkenyl,

- 9) C<sub>2</sub>-C<sub>10</sub> alkynyl,  
 10) heterocyclyl,  
 11) C<sub>3</sub>-C<sub>8</sub> cycloalkyl,  
 12) SO<sub>2</sub>R<sup>a</sup>, and  
 5 13) (C=O)NR<sup>b</sup><sub>2</sub>,

said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one, two or three substituents selected from R<sup>11</sup>, or

- 10 R<sup>12</sup> and R<sup>13</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

- 15 R<sup>a</sup> is (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, aryl, or heterocyclyl;

R<sup>b</sup> is H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, (C=O)OC<sub>1</sub>-C<sub>6</sub> alkyl, (C=O)C<sub>1</sub>-C<sub>6</sub> alkyl or S(O)<sub>2</sub>R<sup>a</sup>;

- 20 R<sup>c</sup> and R<sup>c'</sup> are independently selected from: H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl and (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl; or

- 25 R<sup>c</sup> and R<sup>c'</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

- 30 R<sup>d</sup> and R<sup>d'</sup> are independently selected from: (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>1</sub>-C<sub>6</sub>)alkoxy and NR<sup>b</sup><sub>2</sub>, or

R<sup>d</sup> and R<sup>d'</sup> can be taken together with the phosphorous to which they are attached to form a monocyclic heterocycle with 5-7 members the ring and optionally containing, in addition to the phosphorous, one or two additional heteroatoms selected from NR<sup>e</sup>,

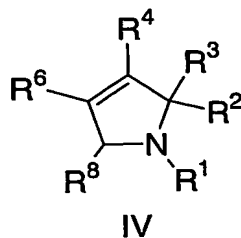


O and S, said monocyclic heterocycle optionally substituted with one, two or three substituents selected from  $R^{11}$ ; and

$R^e$  is selected from: H and  $(C_1-C_6)$ alkyl.

5

6. The compound according to Claim 4 of the Formula IV,



or a pharmaceutically acceptable salt or stereoisomer thereof, wherein

- 10 a is 0 or 1;  
 b is 0 or 1;  
 m is 0, 1, or 2;  
 r is 0 or 1;  
 s is 0 or 1;

15

$R^1$  is selected from:

- 1)  $(C=O)C_1-C_{10}$  alkyl,
- 2)  $(C=O)$ aryl,
- 3)  $(C=O)C_3-C_8$  cycloalkyl,
- 20 4)  $(C=O)$ heterocyclyl,
- 5)  $(C=O)NR^cR^c'$ ,
- 6)  $(C=S)NR^cR^c'$ ,
- 7)  $SO_2NR^cR^c'$ ,
- 8)  $SO_2C_1-C_{10}$  alkyl,
- 25 9)  $SO_2$ -aryl, and
- 10)  $SO_2$ -heterocyclyl,

said alkyl, aryl, cycloalkyl, and heterocyclyl is optionally substituted with one or more substituents selected from  $R^{10}$ ;

R<sup>2</sup> is selected from:

- 1) aryl,
- 2) C<sub>1</sub>-C<sub>6</sub> aralkyl,
- 3) C<sub>3</sub>-C<sub>8</sub> cycloalkyl, and
- 5 4) heterocyclyl,

said aryl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>3</sup>, R<sup>4</sup> and R<sup>8</sup> are independently selected from:

- 10 1) H,
- 2) C<sub>1</sub>-C<sub>10</sub> alkyl, and
- 3) C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

15

R<sup>6</sup> is selected from:

- 1) aryl,
- 2) C<sub>1</sub>-C<sub>6</sub> aralkyl,
- 3) C<sub>3</sub>-C<sub>8</sub> cycloalkyl, and
- 20 4) heterocyclyl,

said aryl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>10</sup> is independently selected from:

- 25 1) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 2) (C=O)<sub>a</sub>O<sub>b</sub>aryl,
- 3) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 4) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 5) (C=O)<sub>a</sub>O<sub>b</sub> heterocyclyl,
- 30 6) CO<sub>2</sub>H,
- 7) halo,
- 8) CN,
- 9) OH,
- 10) O<sub>b</sub>C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,

- 11)  $O_a(C=O)_bNR^{12}R^{13}$ ,  
 12)  $S(O)_mR^a$ ,  
 13)  $S(O)_2NR^{12}R^{13}$ ,  
 14) oxo,  
 5 15) CHO,  
 16)  $(N=O)R^{12}R^{13}$ , or  
 17)  $(C=O)_aO_bC_3-C_8$  cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one, two or three substituents selected from  $R^{11}$ ;

10

$R^{11}$  is selected from:

- 1)  $(C=O)_rO_s(C_1-C_{10})$ alkyl,  
 2)  $O_r(C_1-C_3)$ perfluoroalkyl,  
 3) oxo,  
 15 4) OH,  
 5) halo,  
 6) CN,  
 7)  $(C_2-C_{10})$ alkenyl,  
 8)  $(C_2-C_{10})$ alkynyl,  
 20 9)  $(C=O)_rO_s(C_3-C_6)$ cycloalkyl,  
 10)  $(C=O)_rO_s(C_0-C_6)$ alkylene-aryl,  
 11)  $(C=O)_rO_s(C_0-C_6)$ alkylene-heterocyclyl,  
 12)  $(C=O)_rO_s(C_0-C_6)$ alkylene- $N(R^b)_2$ ,  
 13)  $C(O)R^a$ ,  
 25 14)  $(C_0-C_6)$ alkylene- $CO_2R^a$ ,  
 15)  $C(O)H$ ,  
 16)  $(C_0-C_6)$ alkylene- $CO_2H$ ,  
 17)  $C(O)N(R^b)_2$ ,  
 18)  $S(O)_mR^a$ , and  
 30 19)  $S(O)_2N(R^b)_2$ ;

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, and heterocyclyl is optionally substituted with up to three substituents selected from  $R^b$ , OH,  $(C_1-C_6)$ alkoxy, halogen,  $CO_2H$ , CN,  $O(C=O)C_1-C_6$  alkyl, oxo, and  $N(R^b)_2$ ;

R<sup>12</sup> and R<sup>13</sup> are independently selected from:

- 1) H,
- 2) (C=O)O<sup>b</sup>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 5 3) (C=O)O<sup>b</sup>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 4) (C=O)O<sup>b</sup>aryl,
- 5) (C=O)O<sup>b</sup>heterocyclyl,
- 6) C<sub>1</sub>-C<sub>10</sub> alkyl,
- 7) aryl,
- 10 8) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 9) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 10) heterocyclyl,
- 11) C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 12) SO<sub>2</sub>R<sup>a</sup>, and
- 15 13) (C=O)NR<sup>b</sup><sub>2</sub>,

said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one, two or three substituents selected from R<sup>11</sup>, or

20 R<sup>12</sup> and R<sup>13</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

25 R<sup>a</sup> is independently selected from: (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, aryl, and heterocyclyl;

R<sup>b</sup> is independently selected from: H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, (C=O)OC<sub>1</sub>-C<sub>6</sub> alkyl, (C=O)C<sub>1</sub>-C<sub>6</sub> alkyl or S(O)<sub>2</sub>R<sup>a</sup>; and

30 R<sup>c</sup> and R<sup>c'</sup> are independently selected from: H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl and (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl or

$R^C$  and  $R^{C'}$  can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from  $R^{11}$ .

7. The compound according to Claim 6 or a pharmaceutically acceptable salt or stereoisomer thereof, wherein:

10  $R^1$  is selected from:

- 1)  $(C=O)NR^C R^{C'}$ ,
- 2)  $SO_2NR^C R^{C'}$ , and
- 3)  $SO_2C_1-C_{10}$  alkyl,

15 said alkyl, is optionally substituted with one, two or three substituents selected from  $R^{10}$ ;

$R^2$  is selected from:

- 1) aryl, and
- 2) heteroaryl,

20 said aryl and heteroaryl is optionally substituted with one or more substituents selected from  $R^{10}$ ;

$R^3$ ,  $R^4$  and  $R^8$  are independently selected from:

- 1) H, and
- 25 2)  $C_1-C_{10}$  alkyl,

said alkyl is optionally substituted with one or more substituents selected from  $R^{10}$ ;

$R^6$  is selected from:

- 1) aryl, and
- 30 2) heterocyclyl,

said alkyl, aryl and heterocyclyl is optionally substituted with one or more substituents selected from  $R^{10}$ ; and

$R^{10}$ ,  $R^{11}$ ,  $R^{12}$ ,  $R^{13}$ ,  $R^a$ ,  $R^b$ ,  $R^c$  and  $R^{c'}$  are as described immediately above.

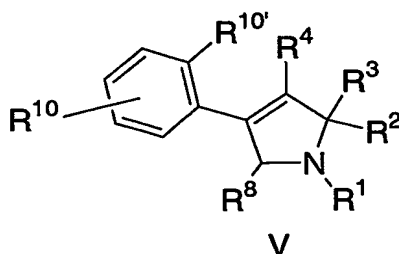
8. The compound according to Claim 7 or a pharmaceutically acceptable salt or stereoisomer thereof, wherein:

- 5 R<sup>2</sup> and R<sup>6</sup> are independently selected from phenyl or pyridyl, optionally substituted with one or two substituents selected from R<sup>10</sup>; and  
R<sup>1</sup>, R<sup>3</sup>, R<sup>4</sup> and R<sup>8</sup> are as described in Claim 7.

9. The compound according to Claim 7, or the pharmaceutically acceptable salt or stereoisomer thereof, wherein R<sup>2</sup> is phenyl, optionally substituted with one or two substituents selected from R<sup>10</sup>; and

R<sup>1</sup>, R<sup>3</sup>, R<sup>4</sup>, R<sup>6</sup> and R<sup>8</sup> are as described in Claim 7.

10. The compound according to Claim 5 of the Formula V,



wherein:

- a is 0 or 1;  
20 b is 0 or 1;  
m is 0, 1, or 2;  
r is 0 or 1;  
s is 0 or 1;

- 25 R<sup>1</sup> is selected from:

- 1) (C=O)C<sub>1</sub>-C<sub>10</sub> alkyl,
- 2) (C=O)aryl,
- 3) (C=O)C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 4) (C=O)heterocyclyl,

- 5)  $(C=O)NR^cR^{c'}$ ,
- 6)  $(C=S)NR^cR^{c'}$ ,
- 7)  $SO_2NR^cR^{c'}$ ,
- 8)  $SO_2C_1-C_{10}$  alkyl,
- 5 9)  $SO_2$ -aryl, and
- 10)  $SO_2$ -heterocyclyl,

said alkyl, aryl, cycloalkyl, and heterocyclyl is optionally substituted with one or more substituents selected from  $R^{10}$ ;

10  $R^2$  is selected from:

- 1) aryl,
- 2)  $C_1-C_6$  aralkyl,
- 3)  $C_3-C_8$  cycloalkyl, and
- 4) heterocyclyl,

15 said aryl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from  $R^{10}$ ;

$R^3$ ,  $R^4$  and  $R^8$  are independently selected from:

- 1) H,
- 20 2)  $C_1-C_{10}$  alkyl, and
- 3)  $C_1-C_6$  perfluoroalkyl,

said alkyl is optionally substituted with one or more substituents selected from  $R^{10}$ ;

$R^{10}$  is independently selected from:

- 25 1)  $(C=O)_aO_bC_1-C_{10}$  alkyl,
- 2)  $(C=O)_aO_b$ aryl,
- 3)  $C_2-C_{10}$  alkenyl,
- 4)  $C_2-C_{10}$  alkynyl,
- 5)  $(C=O)_aO_b$  heterocyclyl,
- 30 6)  $CO_2H$ ,
- 7) halo,
- 8) CN,
- 9) OH,
- 10)  $O_bC_1-C_6$  perfluoroalkyl,

- 11)  $O_a(C=O)_bNR^{12}R^{13}$ ,  
 12)  $S(O)_mR^a$ ,  
 13)  $S(O)_2NR^{12}R^{13}$ ,  
 14) oxo,  
 5 15) CHO,  
 16)  $(N=O)R^{12}R^{13}$ , or  
 17)  $(C=O)_aO_bC_3-C_8$  cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one, two or three substituents selected from  $R^{11}$ ;

10  $R^{10'}$  is halogen;

$R^{11}$  is selected from:

- 1)  $(C=O)_rO_s(C_1-C_{10})$ alkyl,  
 2)  $O_r(C_1-C_3)$ perfluoroalkyl,  
 15 3) oxo,  
 4) OH,  
 5) halo,  
 6) CN,  
 7)  $(C_2-C_{10})$ alkenyl,  
 20 8)  $(C_2-C_{10})$ alkynyl,  
 9)  $(C=O)_rO_s(C_3-C_6)$ cycloalkyl,  
 10)  $(C=O)_rO_s(C_0-C_6)$ alkylene-aryl,  
 11)  $(C=O)_rO_s(C_0-C_6)$ alkylene-heterocyclyl,  
 12)  $(C=O)_rO_s(C_0-C_6)$ alkylene- $N(R^b)_2$ ,  
 25 13)  $C(O)R^a$ ,  
 14)  $(C_0-C_6)$ alkylene- $CO_2R^a$ ,  
 15)  $C(O)H$ ,  
 16)  $(C_0-C_6)$ alkylene- $CO_2H$ ,  
 17)  $C(O)N(R^b)_2$ ,  
 30 18)  $S(O)_mR^a$ , and  
 19)  $S(O)_2N(R^b)_2$ ;

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, alkylene and heterocyclyl is optionally substituted with up to three substituents selected from  $R^b$ , OH,  $(C_1-C_6)$ alkoxy, halogen,  $CO_2H$ , CN,  $O(C=O)C_1-C_6$  alkyl, oxo, and  $N(R^b)_2$ ;



R<sup>12</sup> and R<sup>13</sup> are independently selected from:

- 1) H,
- 2) (C=O)O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 5 3) (C=O)O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 4) (C=O)O<sub>b</sub>aryl,
- 5) (C=O)O<sub>b</sub>heterocyclyl,
- 6) C<sub>1</sub>-C<sub>10</sub> alkyl,
- 7) aryl,
- 10 8) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 9) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 10) heterocyclyl,
- 11) C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 12) SO<sub>2</sub>R<sup>a</sup>, and
- 15 13) (C=O)NR<sup>b</sup><sub>2</sub>,

said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one, two or three substituents selected from R<sup>11</sup>, or

20 R<sup>12</sup> and R<sup>13</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

25 R<sup>a</sup> is independently selected from: (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, aryl, and heterocyclyl;

R<sup>b</sup> is independently selected from: H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, (C=O)OC<sub>1</sub>-C<sub>6</sub> alkyl, (C=O)C<sub>1</sub>-C<sub>6</sub> alkyl or S(O)<sub>2</sub>R<sup>a</sup>; and

30 R<sup>c</sup> and R<sup>c'</sup> are independently selected from: H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl and (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl or

RC and RC' can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with  
5 one, two or three substituents selected from R<sup>11</sup>.

11. A compound selected from:

- 10 4-(2-chloro-5-fluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- (+)-4-(2,5-difluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 15 (-)-4-(2,5-difluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 4-(5-chloro-2-fluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 20 4-(2-fluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 4-(2-fluoro-5-methylphenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 25 4-(5-bromo-2-fluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 4-[[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]morpholine;
- 30 4-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]morpholine;
- N,N-dimethyl-2,4-diphenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

- 3-[2-fluoro-5-(trifluoromethyl)phenyl]-N,N-dimethyl-5-phenyl-2,3-dihydro-1H-pyrrole-1-carboxamide;
- 5 2-(3-fluorophenyl)-4-(2,5-difluorophenyl)- N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 4-(2,5-difluorophenyl)-2-(4-fluorophenyl)-N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 10 4-(2,5-difluorophenyl)-2-(2-fluorophenyl)-N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 2-(3-bromophenyl)-4-(2,5-difluorophenyl)- N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 15 2-(3-aminophenyl)-4-(2,5-difluorophenyl)- N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 20 4-(2,5-difluorophenyl)-2-(3-methylphenyl)-N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- (2S)-4-(5-chloro-2-fluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 25 4-(2,5-difluorophenyl)-1-(methylsulfonyl)-2-phenyl-2,5-dihydro-1H-pyrrole;
- 4-(2,5-difluorophenyl)-1-(ethylsulfonyl)-2-phenyl-2,5-dihydro-1H-pyrrole;
- 30 4-(2,5-difluorophenyl)-2-phenyl-1-(propylsulfonyl)-2,5-dihydro-1H-pyrrole;
- 4-(2,5-difluorophenyl)-1-(isopropylsulfonyl)-2-phenyl-2,5-dihydro-1H-pyrrole;
- 4-(5-chloro-2-fluorophenyl)-1-(methylsulfonyl)-2-phenyl-2,5-dihydro-1H-pyrrole;
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- 4-(5-chloro-2-fluorophenyl)-1-(isopropylsulfonyl)-2-phenyl-2,5-dihydro-1H-pyrrole;
- 4-(2-fluoro-5-methylphenyl)-1-(isopropylsulfonyl)-2-phenyl-2,5-dihydro-1H-pyrrole;
- 5 2-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]sulfonyl}ethanamine;
- 2-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]sulfonyl}-N,N-dimethylethanamine;
- 10 1-acetyl-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrole;
- 4-(2-chloro-5-fluorophenyl)-1-pivaloyl-2-phenyl-2,5-dihydro-1H-pyrrole;
- 15 4-(2-chloro-5-fluorophenyl)-1-isobutyryl-2-phenyl-2,5-dihydro-1H-pyrrole;
- 4-(2,5-difluorophenyl)-1-(2,2-dimethylpropanoyl)-2-phenyl-2,5-dihydro-1H-pyrrole;
- 1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-methyl-1-
- 20 oxopropan-2-ol;
- 1-[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-methyl-1-
- oxopropan-2-ol;
- 25 1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-methyl-1-
- oxopropan-2-amine;
- 4-(2-fluoro-5-isocyanophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-
- carboxamide;
- 30 4-(2,5-difluorophenyl)-2-phenyl-1-(trifluoroacetyl)-2,5-dihydro-1H-pyrrole;
- 4-(5-chloro-2-fluorophenyl)-2-phenyl-1-(trifluoroacetyl)-2,5-dihydro-1H-pyrrole;

(1S)-1-{{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-2-methylpropylamine;

5 (1R)-1-{{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-2-methylpropylamine;

4-(2,5-difluorophenyl)-2-phenyl-1-L-prolyl-2,5-dihydro-1H-pyrrole;

10 4-(2,5-difluorophenyl)-2-phenyl-1-D-prolyl-2,5-dihydro-1H-pyrrole;

(4R)-4-{{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-1,3-thiazolidine;

15 methyl (3S)-3-amino-4-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-4-oxobutanoate;

(4S)-4-amino-5-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-5-oxopentanamide;

20 (1S)-1-{{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-3-(methylthio)propylamine;

(1S)-1-{{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-3-(methylsulfonyl)propylamine;

25 (2S)-2-{{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}piperidine;

30 (1S)-1-{{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}pentylamine;

(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-(thien-2-ylmethyl)ethylamine;

- 4-{{4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl}carbonyl}-1,1-dioxidotetrahydro-2H-thiopyran-4-ylamine;
- (2S)-1-{{4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl}-N-methyl-1-oxopropan-2-amine;
- (1S)-1-{{4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl}carbonyl}propylamine;
- (1S)-2-{{4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl}-2-oxo-1-phenylethanamine;
- (1S)-2-{{4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl}-2-oxo-1-phenylethanamine;
- (4S)-4-amino-5-{{4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl}-5-oxopentanamide  
3-{{4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl}-3-oxopropan-1-amine;
- (1S,2S)-1-{{4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl}carbonyl}-2-methylbutylamine;
- (1S)-1-{{4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl}carbonyl}butylamine;
- (1S)-1-cyclopropyl-2-{{4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl}-2-oxoethanamine;
- 1-{{4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl}carbonyl}cyclopropanamine;
- 1-{{4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl}-1-oxopropan-2-amine;

- (1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-methyl-2-oxoethylamine;
- 5 (1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-(pyridin-2-ylmethyl)ethylamine;
- (1S)-1-cyclohexyl-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;
- 10 (1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-(4-iodobenzyl)-2-oxoethylamine;
- (1S)-1-benzyl-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethylamine;
- 15 4-{(2S)-2-amino-3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxopropyl}phenol;
- (3S)-3-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-1,2,3,4-tetrahydroisoquinoline;
- 20 (1S)-1-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-3-phenylpropylamine;
- (1S)-1-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-3-methylbutylamine;
- 25 (1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-(pyridin-3-ylmethyl)ethylamine;
- 30 1-[(2S)-azetidin-2-ylcarbonyl]-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrole;
- (3S)-3-amino-4-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-4-oxobutanamide;
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4-(2,5-difluorophenyl)-1-[(2S)-2,5-dihydro-1H-pyrrol-2-ylcarbonyl]-2-phenyl-2,5-dihydro-1H-pyrrole;

5 4-(2,5-difluorophenyl)-1-[(2-methylazetidin-2-yl)carbonyl]-2-phenyl-2,5-dihydro-1H-pyrrole;

(1S)-1-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-2,2-dimethylpropylamine;  
methyl (4S)-4-amino-5-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-5-oxopentanoate;

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4-(2,5-difluorophenyl)-2-phenyl-1-[(2S,3S)-2-phenylpyrrolidin-3-yl]carbonyl}-2,5-dihydro-1H-pyrrole;

15 4-(2,5-difluorophenyl)-2-phenyl-1-[(5-phenylpyrrolidin-3-yl)carbonyl]-2,5-dihydro-1H-pyrrole;

(2S)-2-amino-3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxopropan-1-ol;

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(2R,3S)-3-amino-4-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-4-oxobutan-2-ol;

(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-(methoxymethyl)-2-oxoethylamine;

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4-(2,5-difluorophenyl)-2-phenyl-1-(pyrrolidin-3-ylcarbonyl)-2,5-dihydro-1H-pyrrole;

4-(2,5-difluorophenyl)-2-phenyl-1-[(3-phenylpyrrolidin-3-yl)acetyl]-2,5-dihydro-1H-pyrrole;

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(1S)-1-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-3,3-difluoropropylamine;



- (1S)-3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxo-1-phenylpropan-1-amine;
- 4-(2,5-difluorophenyl)-2-phenyl-1-[(4S)-4-phenyl-L-prolyl]-2,5-dihydro-1H-pyrrole;
- 5 1-{2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}cyclohexanamine;
- 2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;
- 10 4-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}piperidin-4-amine;
- (1S,3R)-3-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}cyclopentanamine;
- 15 (1R,4S)-4-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}cyclopent-2-en-1-amine;
- (1S,4R)-4-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}cyclopent-2-en-1-amine;
- 20 (1S)-1-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}but-3-ynylamine;
- (1R)-3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxo-1-phenylpropan-1-amine;
- 25 3-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-2-phenylpiperidine;
- 30 (1S)-1-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}but-3-enylamine;
- (2S)-3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-(methylamino)-3-oxopropan-1-ol;
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(3R,5S)-5-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}pyrrolidin-3-ol;

5 (1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-(1,3-thiazol-4-ylmethyl)ethylamine;

(1R)-1-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}but-3-enylamine;

10 (2S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,3-dimethyl-1-oxobutan-2-amine;

(2S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,4-dimethyl-1-oxopentan-2-amine;

15 (1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-[(1-methyl-1H-imidazol-4-yl)methyl]-2-oxoethylamine;

4-(2,5-difluorophenyl)-1-(N~6~-formyl-L-lysyl)-2-phenyl-2,5-dihydro-1H-pyrrole;

20 (2S,3S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,3-dimethyl-1-oxopentan-2-amine;

(1S)-1-(cyclohexylmethyl)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethylamine;

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(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-(1H-indol-3-ylmethyl)-2-oxoethylamine;

30 (1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-(isocyanomethyl)-2-oxoethylamine;

(1S)-1-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-3,3-dimethylbutylamine;

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- 1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,3-dimethyl-1-oxobutan-2-amine;
- 5 1-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}cyclohexanamine;
- 1-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}cyclopentanamine;
- 10 (1S)-3-(benzyloxy)-1-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}propylamine;
- 1-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,3-dimethyl-1-oxobutan-2-amine;
- 15 1-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}cyclopent-3-en-1-amine;
- (1S)-1-cyclopentyl-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;
- 20 4-(2,5-difluorophenyl)-1-(2-methylpropyl)-2-phenyl-2,5-dihydro-1H-pyrrole;
- 1-4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-methyl-1-oxopropan-2-amine;
- 25 (1S)-1-{[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-2-methylpropylamine;
- (1S)-2-[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-cyclopropyl-2-oxoethanamine;
- 30 (1S,2S)-1-{[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-2-methylbutylamine;
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- (1S)-1-{[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}pentylamine;
- 5 (1S)-1-{[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-3,3-dimethylbutylamine;
- (1S)-1-{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-2,2-dimethylpropylamine;
- 10 (1S)-1-{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-2-methylpropylamine;
- (1S)-1-cyclohexyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;
- 15 (1S)-1-{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}but-3-enylamine;
- (1S)-1-{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}but-3-ynylamine;
- 20 (1S)-1-benzyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethylamine;
- (1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;
- 25 1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-methyl-1-oxopropan-2-amine;
- 30 (1S)-1-{[(2S)-4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-2,2-dimethylpropylamine;
- (1S)-1-{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}pentylamine;
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- (1S)-1-[[ (2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-3-methylbutylamine;
- 5 (1S)-1-[[ (2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-3,3-dimethylbutylamine;
- 1-cyclopropyl-3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxopropan-1-amine;
- 10 (1S)-2-[(2S)-4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-cyclopropyl-2-oxoethanamine;
- (1S)-1-[[ (2S)-4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2-methylpropylamine;
- 15 (1S,2S)-1-[[ (2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2-methylbutylamine;
- 4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-1-(2-methylalanyl)-2,5-dihydro-1H-pyrrole;
- 20 (2S)-4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-1-(2-methylalanyl)-2,5-dihydro-1H-pyrrole;
- 25 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-1-L-valyl-2,5-dihydro-1H-pyrrole;
- 4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-1-L-valyl-2,5-dihydro-1H-pyrrole;
- (2S)-4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-1-L-valyl-2,5-dihydro-1H-pyrrole;
- 30 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-1-(2-methylalanyl)-2,5-dihydro-1H-pyrrole;

- 3-[1-[(2S)-2-amino-2-cyclopropylethanoyl]-4-(5-chloro-2-fluorophenyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol;
- 5 4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-1-L-isoleucyl-2,5-dihydro-1H-pyrrole;
- 4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-1-L-norleucyl-2,5-dihydro-1H-pyrrole;
- 10 (2S)-4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-1-(3-methyl-L-valyl)-2,5-dihydro-1H-pyrrole;
- (2S)-4-(2,5-Difluorophenyl)-N-methyl-2-phenyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 15 (2S)-4-(2,5-Difluorophenyl)-N-methyl-2-phenyl-N-(piperidin-4-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;
- (2S)-4-(5-Chloro-2-fluorophenyl)-N-methyl-2-phenyl-N-[(3R)-pyrrolidin-3-yl]-2,5-
- 20 dihydro-1H-pyrrole-1-carboxamide;
- (2S)-4-(5-chloro-2-fluorophenyl)-N-methyl-2-phenyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 25 (2S)-4-(5-chloro-2-fluorophenyl)-N-methyl-2-phenyl-N-[(3S)-pyrrolidin-3-yl]-2,5-dihydro-1H-pyrrole-1-carboxamide;
- (2S)-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-pyrrolidin-3-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 30 (2S)-4-(2,5-difluorophenyl)-N-methyl-N-[(3S)-1-methylpyrrolidin-3-yl]-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- (2S)-4-(2,5-difluorophenyl)-N-methyl-N-[(3R)-1-methylpyrrolidin-3-yl]-2-phenyl-
- 35 2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(2,5-difluorophenyl)-N-methyl-N-[(1-methyl-5-oxopyrrolidin-2-yl)methyl]-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

- 5 (2S)-4-(2,5-difluorophenyl)-N-(1,3-dioxolan-2-ylmethyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-tetrahydrofuran-3-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

- 10 (2S)-N-(1-allylpiperidin-4-yl)-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

allyl 4-[[[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methyl)amino]piperidine-1-carboxylate;

- 15 allyl 4-[[[[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methyl)amino]methyl]piperidine-1-carboxylate;

- 20 (2S)-4-(2,5-difluorophenyl)-N-methyl-N-(1-methylpiperidin-4-yl)-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-Difluorophenyl)-N-methyl-N-[(1-methylpiperidin-3-yl)methyl]-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

- 25 4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-(pyridin-3-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;

N-benzyl-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

- 30 4-(2,5-difluorophenyl)-N-methyl-N-[(1-methyl-1H-pyrazol-4-yl)methyl]-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

- 35 4-(2,5-difluorophenyl)-N-[2-(dimethylamino)ethyl]-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

- 4-(2,5-difluorophenyl)-N-(2-hydroxyethyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 5 4-(2,5-difluorophenyl)-N-isobutyl-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-(2-pyridin-2-ylethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 10 4-(2,5-difluorophenyl)-N-(2-methoxyethyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 4-(2,5-difluorophenyl)-N-(2,3-dihydroxypropyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 15 4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-(2-phenylethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 4-(2,5-difluorophenyl)-2-phenyl-N-propyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 20 4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 25 1-Acetyl-4-(2,5-difluorophenyl)-2-methyl-2-phenyl-2,5-dihydro-1H-pyrrole;
- (2S)-1-[4-(2,5-difluorophenyl)-2-methyl-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-methyl-1-oxobutan-2-amine;
- 30 (2S)-4-(2,5-difluorophenyl)-N,N,2-trimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;



(2S)-4-(5-chloro-2-fluorophenyl)-N,N,2-trimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

5 (2S)-4-(5-chloro-2-fluorophenyl)-2-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(5-chloro-2-fluorophenyl)-N-ethyl-2-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

10 (2S)-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3,3-dimethyl-1-oxobutan-2-ol;

(2S)-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-methyl-1-oxobutan-2-ol;

15 (2S,3S)-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-methyl-1-oxopentan-2-ol;

20 1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-oxohexan-2-ol;

(2S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-oxo-3-phenylpropan-2-ol;

25 (2S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-4-methyl-1-oxopentan-2-ol;

(1S)-1-cyclohexyl-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanol;

30 (2S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3,3-dimethyl-1-oxobutan-2-ol;

35 N-1-[(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl]-N-2,N-2-dimethylglycinamide;

N-1-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-N-2-methylglycinamide;

5 N-1-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}glycinamide;

N-1-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-2-methylalaninamide;

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N-1-{(1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}glycinamide;

15 N-1-{(1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-N-2,N-2-dimethylglycinamide;

N-1-{(1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-N-2,N-2-dimethylglycinamide, N-oxide;

20 N-1-{(1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-2-methylalaninamide;

N-1-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-N-2,N-2-dimethylglycinamide n-oxide;

25 N-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-2-pyrrolidin-1-ylacetamide;

2-azetidin-1-yl-N-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}acetamide;

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N-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-2-morpholin-4-ylacetamide;

35 N-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-2-piperazin-1-ylacetamide;

N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-(4-methylpiperazin-1-yl)acetamide;

- 5 2-azetidin-1-yl-N-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)acetamide;

N-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-pyrrolidin-1-ylacetamide;

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N-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-piperidin-1-ylacetamide;

- 15 N-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-morpholin-4-ylacetamide;

N-1-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N-2-(2-hydroxyethyl)glycinamide;

- 20 N-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-(4-methylpiperazin-1-yl)acetamide;

N-1-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N-2-isopropylglycinamide;

- 25 N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)acetamide;

N-1-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N-2-ethylglycinamide;

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N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-hydroxyacetamide;

- 35 N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)piperazine-1-carboxamide;

- N-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-N'-piperidin-4-ylurea;
- 5 4-amino-N-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}piperidine-1-carboxamide;
- N-(2-aminoethyl)-N'-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}urea;
- 10 N-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-N'-(3-morpholin-4-ylpropyl)urea;
- N-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-N'-[2-(dimethylamino)ethyl]urea;
- 15 2-azetidin-1-yl-N-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}ethanesulfonamide
- 20 N-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-2-(isopropylamino)ethanesulfonamide;
- N-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-2-pyrrolidin-1-ylethanesulfonamide;
- 25 N-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-2-morpholin-4-ylethanesulfonamide;
- N-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-2-piperazin-1-ylethanesulfonamide;
- 30 N-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-2-(4-methylpiperazin-1-yl)ethanesulfonamide;

N-(tert-butyl)-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetamide;

5 2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-isopropylacetamide;

(2S)-1-(2-azetidin-1-yl-2-oxoethyl)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrole;

10 (2S)-4-(2,5-difluorophenyl)-1-(2-oxo-2-pyrrolidin-1-ylethyl)-2-phenyl-2,5-dihydro-1H-pyrrole;

4-([(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetyl)morpholine;

15 1-([(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetyl)piperazine;

1-([(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetyl)-4-methylpiperazine;

20 2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-isopropylbutanamide;

25 4-{2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]butanoyl}morpholine;

2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-ethylacetamide;

30 N-cyclobutyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetamide;

2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-ethylpropanamide;

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- N-cyclobutyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanamide;
- 2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-methylpropanamide;
- 2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-isopropylpropanamide;
- 10 N-(tert-butyl)-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanamide;
- 4-{2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanoyl}morpholine;
- 15 (3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-ethyl-2,2-dimethyl-4-oxobutanamide;
- (3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,2-dimethyl-4-oxo-N-piperidin-4-ylbutanamide;
- 20 (3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,2-dimethyl-4-oxobutanoic acid;
- (3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,N,2,2-tetramethyl-4-oxobutanamide;
- 25 (1S)-1-{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-2,2-dimethyl-3-oxo-3-piperazin-1-ylpropylamine;
- 30 (3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-isopropyl-2,2-dimethyl-4-oxobutanamide;
- (3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,2,2-trimethyl-4-oxobutanamide;
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(3R)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,N,2,2-tetramethyl-4-oxobutanamide;

5 (3R)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,2-dimethyl-4-oxobutanoic acid;

(1R)-1-{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-2,2-dimethyl-3-oxo-3-piperazin-1-ylpropylamine;

10 2-({(1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}amino)-N-ethylacetamide;

2-({(1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}amino)-N-methylacetamide;

15 2-({(1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}amino)-N,N-dimethylacetamide;

2-({(1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}amino)-N-methyl-N-ethylacetamide;

20 2-({(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}amino)-N-methylacetamide;

25 2-({(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}amino)-N-ethylacetamide;

2-({(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}amino)-N,N-dimethylacetamide;

30 2-({(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}amino)-N-isopropylacetamide;

2-({(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}amino)-N-ethyl-N-methylacetamide;

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- 2-(((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)amino)-N,N-diethylacetamide;
- (1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-N-(2-oxo-2-pyrrolidin-1-ylethyl)ethanamine;
- 5 (1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-(2-morpholin-4-yl-2-oxoethyl)-2-oxoethanamine;
- 10 1-((((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)amino)acetyl)piperidin-4-ol;
- (1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-[2-(4-methylpiperazin-1-yl)-2-oxoethyl]-2-oxoethanamine;
- 15 (1S)-N-(2-azetidin-1-yl-2-oxoethyl)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;
- (1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-[2-(1,1-dioxidothiomorpholin-4-yl)-2-oxoethyl]-2-oxoethanamine;
- 20 (1S)-N-[2-(4-acetylpiperazin-1-yl)-2-oxoethyl]-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;
- (1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-(2-morpholin-4-yl-2-oxoethyl)-2-oxoethanamine;
- 25 (1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-N-(2-oxo-2-pyrrolidin-1-ylethyl)ethanamine;
- 30 2-(((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)amino)-N-isopropylacetamide;
- 2-(dimethylamino)ethyl (1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethylcarbamate;
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- 1-methylpiperidin-4-yl (1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethylcarbamate;
- (2S)-4-cyclopropyl-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 5 (2S)-4-cyclopentyl-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- (1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl 4-methylpiperazine-1-carboxylate;
- 10 1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl 2-morpholin-4-ylethylcarbamate;
- N-[(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}oxy)-carbonyl]glycine;
- 15 (1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl 1-methylpiperidin-4-ylcarbamate;
- (1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethylmethyl(1-methylpiperidin-4-yl)carbamate;
- 20 (1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl 4-dimethylamino)piperidine-1-carboxylate;
- 25 tert-butyl (2S)-4-(2-chloro-5-fluoropyrimidin-4-yl)-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxylate;
- (2S)-4-(5-fluoro-2-methylpyrimidin-4-yl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 30 (2S)-4-(2-chloro-5-fluoropyrimidin-4-yl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

- (2S)-4-(4-chloro-5-methylpyrimidin-2-yl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 5 (2S)-4-(6-chloropyrimidin-4-yl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- (2S)-4-(2-chloropyrimidin-4-yl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 10 (2S)-N,N-dimethyl-4-(4-methylpyridin-3-yl)-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- (2S)-N,N-dimethyl-2-phenyl-4-(1,3-thiazol-2-yl)-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 15 (2S)-N,N-dimethyl-2-phenyl-4-(1,3-thiazol-4-yl)-2,5-dihydro-1H-pyrrole-1-carboxamide;
- (2S)-N,N-dimethyl-2-phenyl-4-(1,2-thiazol-5-yl)-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 20 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- (2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 25 (2S)-4-(2,5-difluorophenyl)-N-(2-hydroxyethyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- (2S)-4-(2,5-difluorophenyl)-N-(2-hydroxyethyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 30 N-([4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl)-N-methyl-beta-alanine;
- methyl N-([4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl)-N-methyl-beta-alaninate;
- 35

- 4-{{4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl}acetyl}morpholin-4-ium;
- 5 3-[(2S)-4-(2,5-difluorophenyl)-1-(2-hydroxy-2-methylpropanoyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol;
- 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-sulfonamide;
- 10 3-[4-(2,5-difluorophenyl)-1-(methylsulfonyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol;
- 3-[4-(2,5-difluorophenyl)-1-(morpholin-4-ylcarbonyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol;
- 15 3-[4-(2,5-difluorophenyl)-1-(2,2-dimethylpropanoyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol;
- (2S)-4-(2,5-difluorophenyl)-1-[(methylsulfonyl)acetyl]-2-phenyl-2,5-dihydro-1H-pyrrole;
- 20 (2S)-4-(2,5-difluorophenyl)-2-phenyl-1-[(phenylsulfonyl)acetyl]-2,5-dihydro-1H-pyrrole;
- 3-[ (2S)-1-[(2S)-2-cyclopropyl-2-hydroxyethanoyl]-4-(2,5-difluorophenyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol;
- 25 3-{ (2S)-4-(2,5-difluorophenyl)-1-[(2S)-2-hydroxy-3,3-dimethylbutanoyl]-2,5-dihydro-1H-pyrrol-2-yl }phenol;
- 30 (1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanol;
- (2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-tetrahydrofuran-3-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 35 (2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-(2-methoxyethyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-tetrahydro-2H-pyran-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

3-[(2S)-4-(2,5-difluorophenyl)-1-(piperidin-1-ylcarbonyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol;

4-(2,5-difluorophenyl)-N-[1-(2-fluoroethyl)piperidin-4-yl]-2-(3-hydroxyphenyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-[[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methyl)amino]piperidinium trifluoroacetate;

2-[[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methyl)amino]ethyl 4-methylpiperazine-1-carboxylate;

3-{4-(2,5-difluorophenyl)-1-[(4-methylpiperazin-1-yl)carbonyl]-2,5-dihydro-1H-pyrrol-2-yl}phenol;

2-[[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methyl)amino]ethyl morpholine-4-carboxylate;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(1-methyl-1H-pyrazol-4-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-(isoxazol-5-ylmethyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

2-[[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methyl)amino]ethyl dimethylaminocarboxylate;

2-[[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methyl)amino]ethyl piperidine-1-carboxylate;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(2-oxopyrrolidin-1-yl)ethyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

- 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(5-oxo-4,5-dihydro-1H-1,2,4-triazol-3-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 5 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-(tetrahydro-2H-pyran-4-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-{[5-(methoxymethyl)-1H-pyrazol-3-yl]methyl}-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 10 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-(1,3-thiazol-4-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(4-methyl-1,2,5-oxadiazol-3-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 15 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-(1,3-thiazol-2-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 20 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-(isoxazol-3-ylmethyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(1H-1,2,4-triazol-1-yl)ethyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 25 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(1H-pyrazol-1-yl)ethyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(1-methyl-5-oxopyrrolidin-2-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 30 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-(1-isoxazol-3-ylethyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

- 4-(2,5-difluorophenyl)-N-(1,3-dioxolan-2-ylmethyl)-2-(3-hydroxyphenyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 5 4-(2,5-difluorophenyl)-N-(1,4-dioxan-2-ylmethyl)-2-(3-hydroxyphenyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(5-methyl-1,3,4-oxadiazol-2-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 10 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(methylsulfonyl)ethyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 2-[[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methyl)amino]ethanesulfonic acid;
- 15 2-hydroxyethyl (1S)-1-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-2,2-dimethylpropylcarbamate;
- 3-hydroxypropyl (1S)-1-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-2,2-dimethylpropylcarbamate;
- 20 2-hydroxyethyl {(1S)-1-isopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}carbamate;
- 25 2-hydroxyethyl {(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}carbamate;
- 4-hydroxybutyl (1S)-1-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-2,2-dimethylpropylcarbamate;
- 30 (2S)-4-(2,5-difluorophenyl)-1-[2-(methylsulfonyl)ethyl]-2-phenyl-2,5-dihydro-1H-pyrrole;
- (2S)-4-(2,5-difluorophenyl)-1-[2-(ethylsulfonyl)ethyl]-2-phenyl-2,5-dihydro-1H-pyrrole;
- 35 1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]pentan-3-one;

- 4-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]butan-2-one;
- 5 4-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-3-methylbutan-2-one;
- 2-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-*N,N*-dimethylethanesulfonamide;
- 10 3-{(2*S*)-4-(2,5-difluorophenyl)-1-[2-(methylsulfonyl)ethyl]-2,5-dihydro-1*H*-pyrrol-2-yl}phenol;
- methyl 3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoate;
- 15 (2*S*)-4-(2,5-difluorophenyl)-1-[2-(ethylsulfonyl)propyl]-2-phenyl-2,5-dihydro-1*H*-pyrrole;
- 3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-*N*-methylpropanamide;
- 20 3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-*N,N*-dimethylpropanamide;
- 25 3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-*N,N*,2-trimethylpropanamide;
- 4-{3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoyl}morpholine;
- 30 1-{3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoyl}-4-(methylsulfonyl)piperazine;
- 1-{3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoyl}piperidin-4-ol;
- 35 methyl 3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoate;
- 40 2-({(1*S*)-1-cyclopropyl-2-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-oxoethyl}oxy)-*N*-ethylacetamide ;
- 4-({(1*S*)-1-cyclopropyl-2-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-oxoethoxy}acetyl)morpholine;
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- 2-((1*S*)-1-cyclopropyl-2-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-oxoethoxy)-*N*-(2-hydroxyethyl)acetamide;
- 5 1-((1*S*)-1-cyclopropyl-2-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-oxoethoxy)acetyl)-4-methylpiperazine;
- 1-((1*S*)-1-cyclopropyl-2-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-oxoethoxy)acetyl)piperazine;
- 10 2-((1*S*)-1-cyclopropyl-2-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-oxoethoxy)-*N*-piperidin-4-ylacetamide;
- 1-((1*S*)-1-cyclopropyl-2-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-oxoethoxy)acetyl)piperidin-4-amine;
- 15 *N*-((1*S*)-1-cyclopropyl-2-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-oxoethyl)-3-morpholin-4-yl-3-oxopropan-1-amine;
- 20 *N*<sup>3</sup>-((1*S*)-1-cyclopropyl-2-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-oxoethyl)-*N*<sup>1</sup>,*N*<sup>1</sup>-dimethyl-β-alaninamide; and
- ((1*S*)-1-((2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl)carbonyl)-2,2-dimethylpropyl)(3-morpholin-4-yl-3-oxopropyl)amine;
- 25 or a pharmaceutically acceptable salt or stereoisomer thereof.

12. The compound according to Claim 11 which is selected from:

- 30 (-)-4-(2,5-difluorophenyl)-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;
- 4-(5-chloro-2-fluorophenyl)-*N,N*-dimethyl-2-phenyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;
- 35 4-(2,5-difluorophenyl)-1-(isopropylsulfonyl)-2-phenyl-2,5-dihydro-1*H*-pyrrole;
- 1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-methyl-1-oxopropan-2-ol;



1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-methyl-1-oxopropan-2-amine;

(1S)-1-{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-  
5 2,2-dimethylpropylamine; and

(2S)-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

10 or a pharmaceutically acceptable salt or stereoisomer thereof.

13. The compound according to Claim 11 which is selected from:

(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-  
15 1-yl]-2-oxoethanamine;

(2S)-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

20 (1S)-1-{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-2,2-dimethylpropylamine;

2-({(1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}amino)-N-ethylacetamide ;

25 (2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(2,5-difluorophenyl)-N-methyl-N-(1-methylpiperidin-4-yl)-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide; and

30 (2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-1-L-valyl-2,5-dihydro-1H-pyrrole

or a pharmaceutically acceptable salt or stereoisomer thereof.

35

## 14. A compound which is:

- (1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine
- 5 (2S)-4-(2,5-difluorophenyl)-1-[(methylsulfonyl)acetyl]-2-phenyl-2,5-dihydro-1H-pyrrole
- 10 (2S)-4-(2,5-difluorophenyl)-2-phenyl-1-[(phenylsulfonyl)acetyl]-2,5-dihydro-1H-pyrrole
- 3-[(2S)-1-[(2S)-2-cyclopropyl-2-hydroxyethanoyl]-4-(2,5-difluorophenyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol
- 15 3-{(2S)-4-(2,5-difluorophenyl)-1-[(2S)-2-hydroxy-3,3-dimethylbutanoyl]-2,5-dihydro-1H-pyrrol-2-yl}phenol
- (1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanol
- 20 (2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-tetrahydrofuran-3-yl-2,5-dihydro-1H-pyrrole-1-carboxamide
- (2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-(2-methoxyethyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide
- 25 (2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-tetrahydro-2H-pyran-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide
- 30 3-[(2S)-4-(2,5-difluorophenyl)-1-(piperidin-1-ylcarbonyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol
- 4-(2,5-difluorophenyl)-N-[1-(2-fluoroethyl)piperidin-4-yl]-2-(3-hydroxyphenyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide
- 35 4-[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methylamino)piperidinium trifluoroacetate

- 2-[[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methylamino)ethyl 4-methylpiperazine-1-carboxylate
- 5 3-{4-(2,5-difluorophenyl)-1-[(4-methylpiperazin-1-yl)carbonyl]-2,5-dihydro-1H-pyrrol-2-yl}phenol
- 2-[[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methylamino)ethyl morpholine-4-carboxylate
- 10 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(1-methyl-1H-pyrazol-4-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide
- 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-(isoxazol-5-ylmethyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide
- 15 2-[[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methylamino)ethyl dimethylaminocarboxylate
- 2-[[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methylamino)ethyl piperidine-1-carboxylate
- 20 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(2-oxopyrrolidin-1-yl)ethyl]-2,5-dihydro-1H-pyrrole-1-carboxamide
- 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(5-oxo-4,5-dihydro-1H-1,2,4-triazol-3-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide
- 25 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-(tetrahydro-2H-pyran-4-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide
- 30 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-{[5-(methoxymethyl)-1H-pyrazol-3-yl)methyl]-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide
- 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-(1,3-thiazol-4-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide
- 35

- 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(4-methyl-1,2,5-oxadiazol-3-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide
- 5 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-(1,3-thiazol-2-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide
- 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-(isoxazol-3-ylmethyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide
- 10 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(1H-1,2,4-triazol-1-yl)ethyl]-2,5-dihydro-1H-pyrrole-1-carboxamide
- 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(1H-pyrazol-1-yl)ethyl]-2,5-dihydro-1H-pyrrole-1-carboxamide
- 15 2,5-dihydro-1H-pyrrole-1-carboxamide
- 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(1-methyl-5-oxopyrrolidin-2-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide
- 20 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-(1-isoxazol-3-ylethyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide
- 4-(2,5-difluorophenyl)-N-(1,3-dioxolan-2-ylmethyl)-2-(3-hydroxyphenyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide
- 25 4-(2,5-difluorophenyl)-N-(1,4-dioxan-2-ylmethyl)-2-(3-hydroxyphenyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide
- 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(5-methyl-1,3,4-oxadiazol-2-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide
- 30 2,5-dihydro-1H-pyrrole-1-carboxamide
- 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(methylsulfonyl)ethyl]-2,5-dihydro-1H-pyrrole-1-carboxamide
- 35 2-[[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methyl)amino]ethanesulfonic acid

- 2-hydroxyethyl (1*S*)-1-{[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl}-2,2-dimethylpropylcarbamate
- 5 3-hydroxypropyl (1*S*)-1-{[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl}-2,2-dimethylpropylcarbamate
- 2-hydroxyethyl {(1*S*)-1-isopropyl-2-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-oxoethyl}carbamate
- 10 2-hydroxyethyl {(1*S*)-1-cyclopropyl-2-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-oxoethyl}carbamate
- 4-hydroxybutyl (1*S*)-1-{[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl}-2,2-dimethylpropylcarbamate
- 15 (2*S*)-4-(2,5-difluorophenyl)-1-[2-(methylsulfonyl)ethyl]-2-phenyl-2,5-dihydro-1*H*-pyrrole
- (2*S*)-4-(2,5-difluorophenyl)-1-[2-(ethylsulfonyl)ethyl]-2-phenyl-2,5-dihydro-1*H*-pyrrole
- 20 1-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]pentan-3-one
- 4-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]butan-2-one
- 25 4-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-3-methylbutan-2-one
- 2-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-*N,N*-dimethylethanesulfonamide
- 30 3-{(2*S*)-4-(2,5-difluorophenyl)-1-[2-(methylsulfonyl)ethyl]-2,5-dihydro-1*H*-pyrrol-2-yl}phenol
- 35 methyl 3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoate
- (2*S*)-4-(2,5-difluorophenyl)-1-[2-(ethylsulfonyl)propyl]-2-phenyl-2,5-dihydro-1*H*-pyrrole
- 40 3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-*N*-methylpropanamide
- 3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-*N,N*-dimethylpropanamide
- 45

- 3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-*N,N*,2-trimethylpropanamide
- 5 4-{3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoyl}morpholine
- 1-{3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoyl}-4-(methylsulfonyl)piperazine
- 10 1-{3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoyl}piperidin-4-ol
- methyl 3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoate
- 15 2-({(1*S*)-1-cyclopropyl-2-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-oxoethyl}oxy)-*N*-ethylacetamide
- 20 4-({(1*S*)-1-cyclopropyl-2-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-oxoethoxy}acetyl)morpholine
- 2-({(1*S*)-1-cyclopropyl-2-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-oxoethoxy}-*N*-(2-hydroxyethyl)acetamide
- 25 1-({(1*S*)-1-cyclopropyl-2-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-oxoethoxy}acetyl)-4-methylpiperazine
- 1-({(1*S*)-1-cyclopropyl-2-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-oxoethoxy}acetyl)piperazine
- 30 2-({(1*S*)-1-cyclopropyl-2-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-oxoethoxy}-*N*-piperidin-4-ylacetamide
- 35 1-({(1*S*)-1-cyclopropyl-2-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-oxoethoxy}acetyl)piperidin-4-amine
- N*-{(1*S*)-1-cyclopropyl-2-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-oxoethyl}-3-morpholin-4-yl-3-oxopropan-1-amine
- 40 *N*<sup>3</sup>-{(1*S*)-1-cyclopropyl-2-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-oxoethyl}-*N*<sup>1</sup>,*N*<sup>1</sup>-dimethyl-β-alaninamide
- ((1*S*)-1-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl)-2,2-dimethylpropyl(3-morpholin-4-yl-3-oxopropyl)amine
- 45

or a pharmaceutically acceptable salt thereof.

15. The compound according to Claim 12 which is the TFA salt of  
5 a compound selected from:

2-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]sulfonyl}-N,N-  
dimethylethanamine;

- 10 1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-methyl-1-  
oxopropan-2-amine;

4-(5-chloro-2-fluorophenyl)-2-phenyl-1-(trifluoroacetyl)-2,5-dihydro-1H-pyrrole;

- 15 (1S)-1-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-2-  
methylpropylamine;

(1R)-1-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-2-  
methylpropylamine;

- 20 4-(2,5-difluorophenyl)-2-phenyl-1-L-prolyl-2,5-dihydro-1H-pyrrole;

4-(2,5-difluorophenyl)-2-phenyl-1-D-prolyl-2,5-dihydro-1H-pyrrole;

- 25 (4R)-4-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-1,3-  
thiazolidine;

methyl (3S)-3-amino-4-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-  
yl]-4-oxobutanoate;

- 30 (4S)-4-amino-5-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-5-  
oxopentanamide;

- 35 (1S)-1-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-3-  
(methylthio)propylamine;

- (1S)-1-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-3-(methylsulfonyl)propylamine;
- 5 (2S)-2-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}piperidine;
- (1S)-1-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}pentylamine;
- 10 (1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-(thien-2-ylmethyl)ethylamine;
- 4-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-1,1-dioxidotetrahydro-2H-thiopyran-4-ylamine;
- 15 (2S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-methyl-1-oxopropan-2-amine;
- 20 (1S)-1-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}propylamine;
- (1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-phenylethanamine;
- 25 (1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-phenylethanamine;
- (4S)-4-amino-5-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-5-oxopentanamide
- 30 3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxopropan-1-amine;
- (1S,2S)-1-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-2-methylbutylamine;
- 35



- (1S)-1-{{4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl}carbonyl}butylamine;
- 5 (1S)-1-cyclopropyl-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;
- 1-{{4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl}carbonyl}cyclopropanamine;
- 10 1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-oxopropan-2-amine;
- (1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-methyl-2-
- 15 oxoethylamine;
- (1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-(pyridin-2-ylmethyl)ethylamine;
- 20 (1S)-1-cyclohexyl-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;
- (1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-(4-iodobenzyl)-2-oxoethylamine;
- 25 (1S)-1-benzyl-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethylamine;
- 4-{{(2S)-2-amino-3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxopropyl}phenol};
- 30 (3S)-3-{{4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl}carbonyl}-1,2,3,4-tetrahydroisoquinoline;
- (1S)-1-{{4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl}carbonyl}-3-
- 35 phenylpropylamine;

- (1S)-1-{{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-3-methylbutylamine;
- 5 (1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-(pyridin-3-ylmethyl)ethylamine;
- 1-[(2S)-azetidin-2-ylcarbonyl]-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrole;
- 10 (3S)-3-amino-4-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-4-oxobutanamide;
- 4-(2,5-difluorophenyl)-1-[(2-methylazetidin-2-yl)carbonyl]-2-phenyl-2,5-dihydro-1H-pyrrole;
- 15 (1S)-1-{{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-2,2-dimethylpropylamine;
- 20 methyl (4S)-4-amino-5-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-5-oxopentanoate;
- 4-(2,5-difluorophenyl)-2-phenyl-1-{{[(2S,3S)-2-phenylpyrrolidin-3-yl]carbonyl}-2,5-dihydro-1H-pyrrole;
- 25 4-(2,5-difluorophenyl)-2-phenyl-1-[(5-phenylpyrrolidin-3-yl)carbonyl]-2,5-dihydro-1H-pyrrole;
- (2S)-2-amino-3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxopropan-1-ol;
- 30 (2R,3S)-3-amino-4-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-4-oxobutan-2-ol;
- (1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-(methoxymethyl)-2-oxoethylamine;
- 35

- 4-(2,5-difluorophenyl)-2-phenyl-1-(pyrrolidin-3-ylcarbonyl)-2,5-dihydro-1H-pyrrole;
- 4-(2,5-difluorophenyl)-2-phenyl-1-[(3-phenylpyrrolidin-3-yl)acetyl]-2,5-dihydro-1H-pyrrole;
- 5 (1S)-1-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-3,3-difluoropropylamine;
- 10 (1S)-3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxo-1-phenylpropan-1-amine;
- 4-(2,5-difluorophenyl)-2-phenyl-1-[(4S)-4-phenyl-L-prolyl]-2,5-dihydro-1H-pyrrole;
- 15 1-{2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}cyclohexanamine;
- 2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;
- 20 4-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}piperidin-4-amine;
- (1S,3R)-3-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}cyclopentanamine;
- 25 (1R,4S)-4-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}cyclopent-2-en-1-amine;
- (1S,4R)-4-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}cyclopent-2-en-1-amine;
- 30 (1S)-1-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}but-3-ynylamine;

- (1R)-3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxo-1-phenylpropan-1-amine;
- 3-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-2-phenylpiperidine;
- 5 (1S)-1-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}but-3-enylamine;
- 10 (2S)-3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-(methylamino)-3-oxopropan-1-ol;
- (3R,5S)-5-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}pyrrolidin-3-ol;
- 15 (1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-(1,3-thiazol-4-ylmethyl)ethylamine;
- (1R)-1-{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}but-3-enylamine;
- 20 (2S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,3-dimethyl-1-oxobutan-2-amine;
- (2S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,4-dimethyl-1-oxopentan-2-amine;
- 25 (1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-[(1-methyl-1H-imidazol-4-yl)methyl]-2-oxoethylamine;
- 30 4-(2,5-difluorophenyl)-1-(N-6-formyl-L-lysyl)-2-phenyl-2,5-dihydro-1H-pyrrole;
- (2S,3S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,3-dimethyl-1-oxopentan-2-amine;
- 35

- (1S)-1-(cyclohexylmethyl)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethylamine;
- 5 (1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-(1H-indol-3-ylmethyl)-2-oxoethylamine;
- (1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-(isocyanomethyl)-2-oxoethylamine;
- 10 (1S)-1-{{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-3,3-dimethylbutylamine;
- 1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,3-dimethyl-1-oxobutan-2-amine;
- 15 1-{{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}cyclohexanamine;
- 1-{{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}cyclopentanamine;
- 20 (1S)-3-(benzyloxy)-1-{{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}propylamine;
- 25 1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,3-dimethyl-1-oxobutan-2-amine;
- 1-{{[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}cyclopent-3-en-1-amine;
- 30 (1S)-1-cyclopentyl-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;
- 4-(2,5-difluorophenyl)-1-(2-methylproyl)-2-phenyl-2,5-dihydro-1H-pyrrole;
- 35

(1S)-2-[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-cyclopropyl-2-oxoethanamine;

5 (1S,2S)-1-{[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-2-methylbutylamine;

(1S)-1-{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}but-3-enylamine;

10 (1S)-1-{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}but-3-ynylamine;

(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;

15 1-cyclopropyl-3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxopropan-1-amine;

20 (1S,2S)-1-{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-2-methylbutylamine;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-1-L-valyl-2,5-dihydro-1H-pyrrole;

25 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-1-(2-methylalanyl)-2,5-dihydro-1H-pyrrole;

3-[1-[(2S)-2-amino-2-cyclopropylethanoyl]-4-(5-chloro-2-fluorophenyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol;

30 4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-1-L-isoleucyl-2,5-dihydro-1H-pyrrole;

(2S)-4-(2,5-Difluorophenyl)-N-methyl-2-phenyl-N-(piperidin-4-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;

35

- (2S)-4-(5-Chloro-2-fluorophenyl)-N-methyl-2-phenyl-N-[(3R)-pyrrolidin-3-yl]-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 5 (2S)-4-(5-chloro-2-fluorophenyl)-N-methyl-2-phenyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- (2S)-4-(5-chloro-2-fluorophenyl)-N-methyl-2-phenyl-N-[(3S)-pyrrolidin-3-yl]-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 10 (2S)-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-pyrrolidin-3-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- (2S)-N-(1-allylpiperidin-4-yl)-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 15 4-(2,5-Difluorophenyl)-N-methyl-N-[(1-methylpiperidin-3-yl)methyl]-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-(pyridin-3-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 20 4-(2,5-difluorophenyl)-N-methyl-N-[(1-methyl-1H-pyrazol-4-yl)methyl]-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 25 4-(2,5-difluorophenyl)-N-[2-(dimethylamino)ethyl]-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-(2-pyridin-2-ylethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 30 4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 35

(2S)-1-[4-(2,5-difluorophenyl)-2-methyl-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-methyl-1-oxobutan-2-amine;

- 5 N-1-[(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl]-N-2,N-2-dimethylglycinamide;

N-1-[(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl]-N-2-methylglycinamide;

10

N-1-[(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl]glycinamide;

- 15 N-1-[(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl]-2-methylalaninamide;

N-[(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl]-2-pyrrolidin-1-ylacetamide;

- 20 2-azetidin-1-yl-N-[(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl]acetamide;

N-[(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl]-2-morpholin-4-ylacetamide;

25

N-[(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl]-2-piperazin-1-ylacetamide;

- 30 N-[(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl]-2-(4-methylpiperazin-1-yl)acetamide;

N-1-[(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl]-N-2-isopropylglycinamide;



N-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}piperazine-1-carboxamide;

5 N-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-N'-piperidin-4-ylurea;

4-amino-N-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}piperidine-1-carboxamide;

10 N-(2-aminoethyl)-N'-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}urea;

N-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-N'-(3-morpholin-4-ylpropyl)urea;

15 N-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-N'-[2-(dimethylamino)ethyl]urea;

20 2-azetidin-1-yl-N-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}ethanesulfonamide

N-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-2-(isopropylamino)ethanesulfonamide;

25 N-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-2-pyrrolidin-1-ylethanesulfonamide;

N-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-2-morpholin-4-ylethanesulfonamide;

30 N-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-2-piperazin-1-ylethanesulfonamide;

35 N-{(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-2-(4-methylpiperazin-1-yl)ethanesulfonamide;

- N-(tert-butyl)-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetamide;
- 5 2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-isopropylacetamide;
- (2S)-1-(2-azetidin-1-yl-2-oxoethyl)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrole;
- 10 (2S)-4-(2,5-difluorophenyl)-1-(2-oxo-2-pyrrolidin-1-ylethyl)-2-phenyl-2,5-dihydro-1H-pyrrole;
- 4-{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetyl}morpholine;
- 15 1-{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetyl}piperazine;
- 1-{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetyl}-4-methylpiperazine;
- 20 2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-isopropylbutanamide;
- 25 4-{2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]butanoyl}morpholine;
- 2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-ethylacetamide;
- 30 N-cyclobutyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetamide;
- 2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-ethylpropanamide;
- 35

- N-cyclobutyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanamide;
- 5 2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-methylpropanamide;
- 2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-isopropylpropanamide;
- 10 N-(tert-butyl)-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanamide;
- 4-{2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanoyl}morpholine;
- 15 (3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-ethyl-2,2-dimethyl-4-oxobutanamide;
- (3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,2-dimethyl-4-oxo-N-piperidin-4-ylbutanamide;
- 20 (3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,2-dimethyl-4-oxobutanoic acid;
- 25 (3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,N,2,2-tetramethyl-4-oxobutanamide;
- (1S)-1-{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-2,2-dimethyl-3-oxo-3-piperazin-1-ylpropylamine;
- 30 (3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-isopropyl-2,2-dimethyl-4-oxobutanamide;

- (3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,2,2-trimethyl-4-oxobutanamide;
- 5 (3R)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,N,2,2-tetramethyl-4-oxobutanamide;
- (3R)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,2-dimethyl-4-oxobutanoic acid;
- 10 (1R)-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl-2,2-dimethyl-3-oxo-3-piperazin-1-ylpropylamine
- (1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl 4-methylpiperazine-1-carboxylate;
- 15 (1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl 1-methylpiperidin-4-ylcarbamate;
- (1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethylmethyl(1-methylpiperidin-4-yl)carbamate;
- 20 (1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl-4-dimethylamino)piperidine-1-carboxylate;
- 25 (2S)-N,N-dimethyl-4-(4-methylpyridin-3-yl)-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 4-[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]acetyl]morpholin-4-ium;
- 30 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 2-[[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methylamino)ethyl-4-methylpiperazine-1-carboxylate;
- 35

- 3-{4-(2,5-difluorophenyl)-1-[(4-methylpiperazin-1-yl)carbonyl]-2,5-dihydro-1H-pyrrol-2-yl}phenol;
- 5 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(1-methyl-1H-pyrazol-4-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 4-(2,5-difluoro-phenyl)-2,5-dihydro-2-(3-hydroxyphenyl)-N-methyl-N--[2-(4-methyl-1-piperazinyl)-2-oxoethyl]-H-pyrrole-1-carboxamide;
- 10 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(5-oxo-4,5-dihydro-1H-1,2,4-triazol-3-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 15 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-{[5-(methoxymethyl)-1H-pyrazol-3-yl]methyl}-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-(1,3-thiazol-4-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 20 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-(1,3-thiazol-2-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(1H-1,2,4-triazol-1-yl)ethyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 25 4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(1H-pyrazol-1-yl)ethyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;
- 4-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]butan-2-one;
- 30 4-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-3-methylbutan-2-one;
- 3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-*N*-methylpropanamide;
- 35 3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-*N,N*-dimethylpropanamide;
- 40 3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-*N,N*,2-trimethylpropanamide;
- 4-{3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoyl}morpholine;
- 45

1-{3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoyl}-4-(methylsulfonyl)piperazine;

5 1-{3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoyl}piperidin-4-ol; and

methyl 3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoate.

10 16. A pharmaceutical composition that is comprised of a compound in accordance with Claim 1 and a pharmaceutically acceptable carrier.

15 17. A method of treating or preventing cancer in a mammal in need of such treatment that is comprised of administering to said mammal a therapeutically effective amount of a compound of Claim 1.

20 18. A method of treating cancer or preventing cancer in accordance with Claim 17 wherein the cancer is selected from cancers of the brain, genitourinary tract, lymphatic system, stomach, larynx and lung.

25 19. A method of treating or preventing cancer in accordance with Claim 17 wherein the cancer is selected from histiocytic lymphoma, lung adenocarcinoma, small cell lung cancers, pancreatic cancer, glioblastomas and breast carcinoma.

20. A process for making a pharmaceutical composition which comprises combining a compound of Claim 1 with a pharmaceutically acceptable carrier.

30 21. The composition of Claim 16 further comprising a second compound selected from:

- 35
- 1) an estrogen receptor modulator,
  - 2) an androgen receptor modulator,
  - 3) a retinoid receptor modulator,
  - 4) a cytotoxic/cytostatic agent,
  - 5) an antiproliferative agent,

- 5                   6)     a prenyl-protein transferase inhibitor,  
                  7)     an HMG-CoA reductase inhibitor,  
                  8)     an HIV protease inhibitor,  
                  9)     a reverse transcriptase inhibitor,  
                  10)    an angiogenesis inhibitor, and  
                  11)    a PPAR- $\gamma$  agonist,  
                  12)    a PPAR- $\delta$  agonists;  
                  13)    an inhibitor of cell proliferation and survival signaling, and  
                  14)    an agent that interferes with a cell cycle checkpoint.

10

22.     The composition of Claim 21, wherein the second compound is an angiogenesis inhibitor selected from the group consisting of a tyrosine kinase inhibitor, an inhibitor of epidermal-derived growth factor, an inhibitor of fibroblast-derived growth factor, an inhibitor of platelet derived growth factor, an MMP  
15 inhibitor, an integrin blocker, interferon- $\alpha$ , interleukin-12, pentosan polysulfate, a cyclooxygenase inhibitor, carboxyamidotriazole, combretastatin A-4, squalamine, 6-O-(chloroacetyl-carbonyl)-fumagillol, thalidomide, angiostatin, troponin-1, and an antibody to VEGF.

20

23.     The composition according to Claim 16 further comprising a proteosome inhibitor.

25

24.     The composition according to Claim 16 further comprising a aurora kinase inhibitor.

25.     The composition according to Claim 16 further comprising a Raf kinase inhibitor.

30

26.     The composition according to Claim 16 further comprising a serine/threonine kinase inhibitor.

27.     The composition according to Claim 16 further comprising an inhibitor of another mitotic kinesin which is not KSP.

28. The composition of Claim 22, wherein the second compound is an estrogen receptor modulator selected from tamoxifen and raloxifene.

29. A method of treating cancer which comprises administering a therapeutically effective amount of a compound of Claim 1 in combination with radiation therapy.

30. A method of treating or preventing cancer that comprises administering a therapeutically effective amount of a compound of Claim 1 in combination with a compound selected from:

- 1) an estrogen receptor modulator,
- 2) an androgen receptor modulator,
- 3) a retinoid receptor modulator,
- 4) a cytotoxic/cytostatic agent,
- 5) an antiproliferative agent,
- 6) a prenyl-protein transferase inhibitor,
- 7) an HMG-CoA reductase inhibitor,
- 8) an HIV protease inhibitor,
- 9) a reverse transcriptase inhibitor,
- 10) an angiogenesis inhibitor,
- 11) PPAR- $\gamma$  agonists,
- 12) PPAR- $\delta$  agonists,
- 13) an inhibitor of inherent multidrug resistance,
- 14) an anti-emetic agent,
- 15) an agent useful in the treatment of anemia,
- 16) an agent useful in the treatment of neutropenia,
- 17) an immunologic-enhancing drug,
- 18) an inhibitor of cell proliferation and survival signaling, and
- 19) an agent that interferes with a cell cycle checkpoint.

31. A method of treating cancer that comprises administering a therapeutically effective amount of a compound of Claim 1 in combination with radiation therapy and a compound selected from:

- 1) an estrogen receptor modulator,
- 2) an androgen receptor modulator,



- 5                   3)     a retinoid receptor modulator,  
                  4)     a cytotoxic/cytostatic agent,  
                  5)     an antiproliferative agent,  
                  6)     a prenyl-protein transferase inhibitor,  
                  7)     an HMG-CoA reductase inhibitor,  
                  8)     an HIV protease inhibitor,  
                  9)     a reverse transcriptase inhibitor,  
                  10)    an angiogenesis inhibitor,  
                  11)    PPAR- $\gamma$  agonists,  
10               12)    PPAR- $\delta$  agonists,  
                  13)    an inhibitor of inherent multidrug resistance,  
                  14)    an anti-emetic agent,  
                  15)    an agent useful in the treatment of anemia,  
                  16)    an agent useful in the treatment of neutropenia,  
15               17)    an immunologic-enhancing drug,  
                  18)    an inhibitor of cell proliferation and survival signaling, and  
                  19)    an agent that interferes with a cell cycle checkpoint.

20               32.     A method of treating or preventing cancer which comprises  
administering a therapeutically effective amount of a compound of Claim 1 and  
paclitaxel or trastuzumab.

25               33.     A method of treating or preventing cancer which comprises  
administering a therapeutically effective amount of a compound of Claim 1 and a  
GPIIb/IIIa antagonist.

                  34.     The method of Claim 33 wherein the GPIIb/IIIa antagonist is  
tirofiban.

30               35.     A method of treating or preventing cancer which comprises  
administering a therapeutically effective amount of a compound of Claim 1 in  
combination with a COX-2 inhibitor.

36. A method of treating or preventing cancer which comprises administering a therapeutically effective amount of a compound of Claim 1 in combination with a proteasome inhibitor.

5 37. A method of treating or preventing cancer which comprises administering a therapeutically effective amount of a compound of Claim 1 in combination with an aurora kinase inhibitor.

10 38. A method of treating or preventing cancer which comprises administering a therapeutically effective amount of a compound of Claim 1 in combination with a Raf kinase inhibitor.

15 39. A method of treating or preventing cancer which comprises administering a therapeutically effective amount of a compound of Claim 1 in combination with a serine/threonine kinase inhibitor.

40. A method of treating or preventing cancer which comprises administering a therapeutically effective amount of a compound of Claim 1 in combination with an inhibitor of a mitotic kinesin that is not KSP.

20 41. A method of modulating mitotic spindle formation which comprises administering a therapeutically effective amount of a compound of Claim 1.

25 42. A method of inhibiting the mitotic kinesin KSP which comprises administering a therapeutically effective amount of a compound of Claim 1.